Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\

Data File : BG050915.D

Acq On : 9 Nov 2021 18:59

Operator : CG/JU Sample : PB140594BS

Misc

ALS Vial : 41 Sample Multiplier: 1

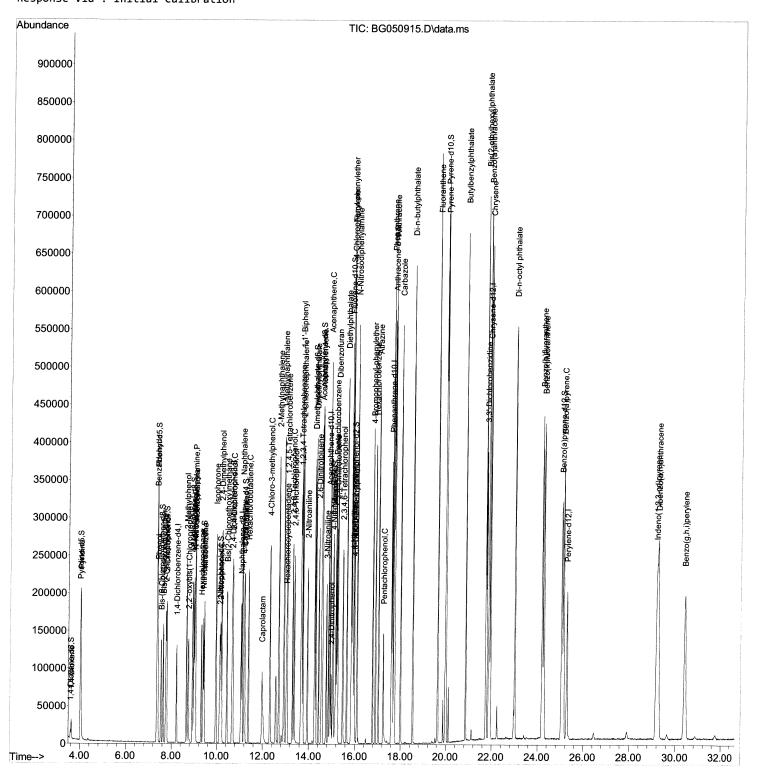
Quant Time: Nov 10 06:59:10 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M

Quant Title : SVOA CALIBRATION

QLast Update : Tue Nov 02 14:49:05 2021 Response via : Initial Calibration Instrument : BNA_G ClientSampleId : SI CS594

Manual IntegrationsAPPROVED



Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\

Data File : BG050915.D

Acq On : 9 Nov 2021 18:59

Operator : CG/JU Sample : PB140594BS

Misc

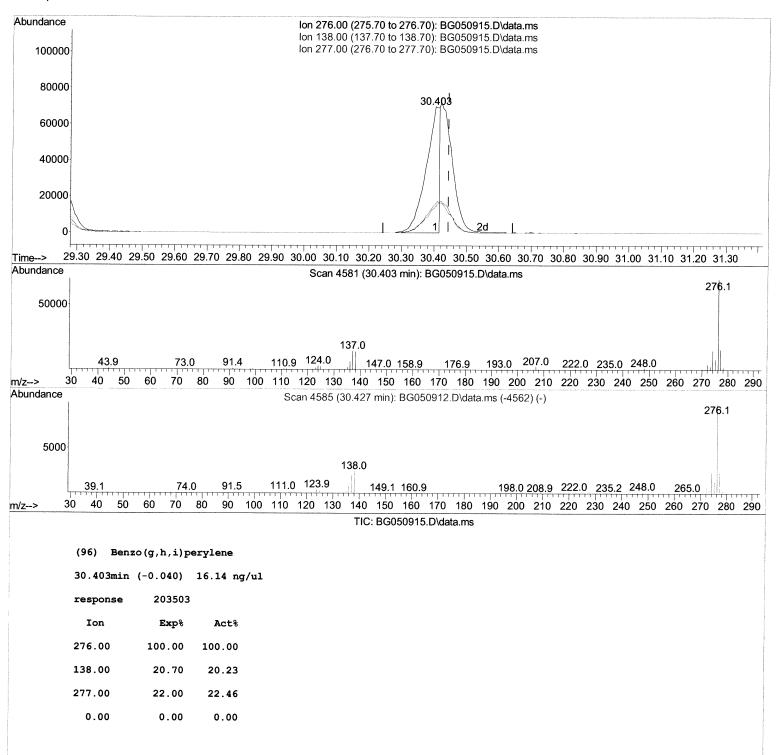
ALS Vial : 41 Sample Multiplier: 1

Quant Time: Nov 10 06:59:10 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M

Quant Title : SVOA CALIBRATION QLast Update : Tue Nov 02 14:49:05 2021 Response via : Initial Calibration Instrument: BNA_G ClientSampleld: SLCS594

Manual IntegrationsAPPROVED



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Operator : CG/JU Sample : PB140594BS

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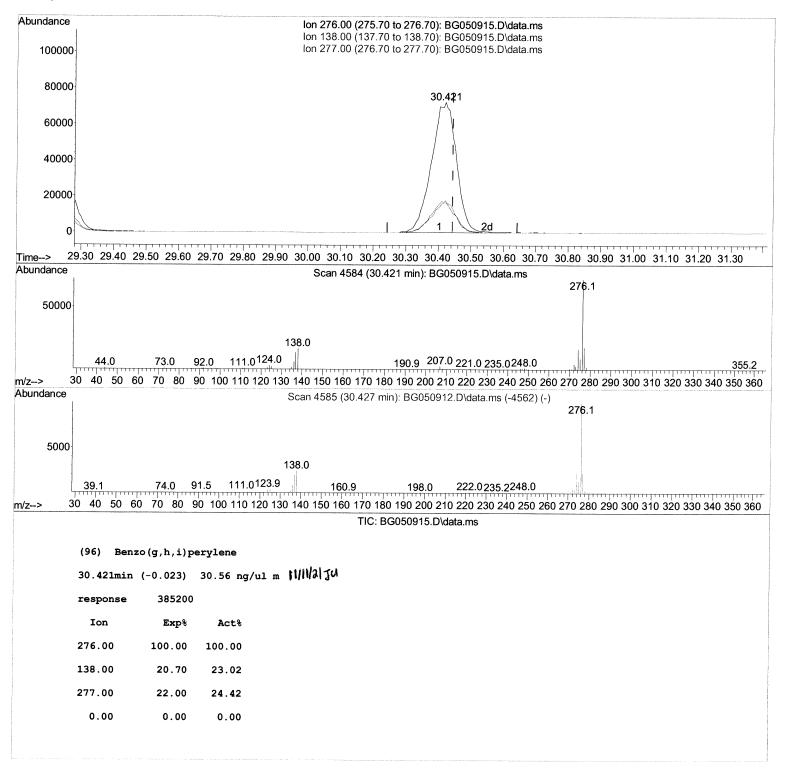
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BNA_G
ClientSampleId:
SLCS594

Manual IntegrationsAPPROVED



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Quant Title : SVOA CALIBRATION QLast Update : Tue Nov 02 14:49:05 2021 Response via : Initial Calibration Instrument : BNA_G ClientSampleId : SLCS594

Manual IntegrationsAPPROVED

Compound	R.T.	QIon	Response	Conc Units De	ev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	8.229	152	35467	20.00 ng/ul	0.00
20) Naphthalene-d8	11.055		158271	20.00 ng/ul	0.00
38) Acenaphthene-d10	14.851		104256	20.00 ng/ul	-0.01
64) Phenanthrene-d10	17.595		231843	20.00 ng/ul	-0.01
79) Chrysene-d12	21.890		196311	20.00 ng/ul	-0.02
88) Perylene-d12	25.292	264	199394	20.00 ng/ul	-0.02
Cusham Manitanian Community					
System Monitoring Compounds	2 502	0.0	6262	·	
3) 1,4-Dioxane-d8	3.593	96	6362	5.79 ng/uL	0.00
4) Pyridine-d57) Phenol-d5	4.011	84	93693	28.50 ng/ul	0.00
9) Bis-(2-Chloroethyl)eth	7.371 7.542	99 67	108459	28.66 ng/ul	0.00
11) 2-Chlorophenol-d4	7.753	67 132	69308	28.36 ng/ul	0.00
15) 4-Methylphenol-d8	8.928	113	76052 83540	29.00 ng/ul 28.05 ng/ul	-0.01 0.00
21) Nitrobenzene-d5	9.404	128	40614	30.20 ng/ul	0.00
24) 2-Nitrophenol-d4	10.127	143	43294	28.95 ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.668	165	75568	30.00 ng/ul	0.00
31) 4-Chloroaniline-d4	11.185	131	95222	24.96 ng/ul	0.00
46) Dimethylphthalate-d6	14.246	166	241050	30.22 ng/ul	0.00
49) Acenaphthylene-d8	14.545	160	301886	30.38 ng/ul	-0.01
54) 4-Nitrophenol-d4	15.039	143	42206	29.18 ng/ul	0.00
60) Fluorene-d10	15.838	176	211847	29.98 ng/ul	-0.01
65) 4,6-Dinitro-2-methylph	15.955	200	38451	27.36 ng/ul	0.00
73) Anthracene-d10	17.695	188	329693	30.08 ng/ul	-0.01
81) Pyrene-d10	19.968	212	379109	29.90 ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.057	264	321380	29.16 ng/ul	-0.02
Target Compounds				0	value
2) 1,4-Dioxane	3.623	88	15473	12.82 ng/uL	96
5) Pyridine	4.034	79	99641	29.28 ng/ul	98
6) Benzaldehyde	7.360	77	76334	31.99 ng/ul	95
8) Phenol	7.401	94	117425	30.00 ng/ul	99
10) Bis(2-Chloroethyl)ether	7.636	93	88790	30.31 ng/ul	99
12) 2-Chlorophenol	7.789	128	80161	30.11 ng/ul	98
13) 2-Methylphenol	8.664	108	85881	29.69 ng/ul	98
<pre>14) 2,2'-oxybis(1-Chloropr</pre>	8.746	45	137132	29.72 ng/ul	98
16) Acetophenone	9.058	105	131923	28.51 ng/ul	99
17) N-Nitroso-di-n-propyla	9.028	70	80955	29.00 ng/ul	98
18) 4-Methylphenol	8.987	108	91487	29.70 ng/ul	97
19) Hexachloroethane	9.322	117	33336	29.95 ng/ul	95
22) Nitrobenzene	9.445	77	118231	31.52 ng/ul	99
23) Isophorone	9.962	82	220342	30.27 ng/ul	99
<pre>25) 2-Nitrophenol</pre>	10.162	139	48448	32.29 ng/ul	99
26) 2,4-Dimethylphenol	10.203	107	103634	31.38 ng/ul	98
<pre>27) Bis(2-Chloroethoxy)met</pre>	10.438	93	121344	30.93 ng/ul	98
29) 2,4-Dichlorophenol	10.691	162	77298	31.46 ng/ul	97
30) Naphthalene	11.102	128	269705	31.16 ng/ul	97
32) 4-Chloroaniline	11.208	127	96545	25.49 ng/ul	99
33) Hexachlorobutadiene	11.373	225	51113	31.69 ng/ul	97
34) Caprolactam	11.966	113	29675	28.45 ng/ul	93
35) 4-Chloro-3-methylphenol	12.313	107	96755	30.82 ng/ul	97

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Misc

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Quant Title : SVOA CALIBRATION QLast Update : Tue Nov 02 14:49:05 2021 Response via : Initial Calibration

Instrument : BNA_G ClientSampleId : SLCS594

Manual IntegrationsAPPROVED

Compound	R.T.	QIon	Response	Conc l	Jnits Dev	(Min)
36) 2-Methylnaphthalene	12.695	142	179549	30.4	 5 ng/ul	100
37) 1-Methylnaphthalene	12.912	142	181018		ng/ul	97
39) 1,2,4,5-Tetrachloroben	13.053	216	96271		ng/ul	95
40) Hexachlorocyclopentadiene	13.029	237	40315		1 ng/ul	96
41) 2,4,6-Trichlorophenol	13.288	196	61723		ng/ul	98
42) 2,4,5-Trichlorophenol	13.364	196	68466		3 ng/ul	99
43) 1,1'-Biphenyl	13.688	154	238654		2 ng/ul	99
44) 2-Chloronaphthalene	13.735	162	186525		3 ng/ul	97
45) 2-Nitroaniline	13.934	65	74445		ng/ul	95
47) Dimethylphthalate	14.293	163	251702		ng/ul	99
48) 2,6-Dinitrotoluene	14.422	165	53041		ng/ul	95
50) Acenaphthylene	14.581	152	312862		ng/ul	99
51) 3-Nitroaniline	14.757	138	50281		ng/ul	92
52) Acenaphthene	14.915	153	207430		ng/ul	96
53) 2,4-Dinitrophenol	14.962	184	21658		ng/ul	91
55) 4-Nitrophenol	15.051	109	39442		ng/ul	98
56) Dibenzofuran	15.250	168	290241		ng/ul	99
57) 2,4-Dinitrotoluene	15.209	165	77159		ng/ul#	97
58) 2,3,4,6-Tetrachlorophenol	15.468	232	52435		ng/ul	96
59) Diethylphthalate	15.644	149	268987		ng/ul	99
61) Fluorene	15.897	166	231737	31.23	ng/ul	99
62) 4-Chlorophenyl-phenyle	15.879	204	121682		ng/ul	98
63) 4-Nitroaniline	15.920	138	54261	31.68	ng/ul	96
66) 4,6-Dinitro-2-methylph	15.973	198	39294	28.67	ng/ul#	96
67) N-Nitrosodiphenylamine	16.091	169	206279	31.83	ng/ul	98
68) 4-Bromophenyl-phenylether	16.778	248	74052	32.11	ng/ul	95
69) Hexachlorobenzene	16.896	284	75723	31.94	ng/ul	98
70) Atrazine	17.031	200	82984	30.20	ng/ul	100
71) Pentachlorophenol	17.242	266	28308		ng/ul	92
72) Phenanthrene	17.642	178	393917		ng/ul	99
74) Anthracene	17.730	178	387998		ng/ul	99
75) 1,2,3,4-Tetrachloroben	13.658	216	99989		ng/uL	97
76) Pentachlorobenzene	15.168	250	90298		ng/uL	98
77) Carbazole	18.000	167	364365		ng/ul	99
78) Di-n-butylphthalate	18.529	149	464130		ng/ul	100
80) Fluoranthene	19.639	202	478638		ng/ul	99
82) Pyrene	19.998	202	463755		ng/ul	99
83) Butylbenzylphthalate	20.861	149	202006		ng/ul	98
84) 3,3'-Dichlorobenzidine	21.778	252	128188		ng/ul	98
85) Benzo(a)anthracene	21.872	228	428101		ng/ul	99
86) Bis(2-ethylhexyl)phtha87) Chrysene	21.743	149	289056		ng/ul	99
	21.943	228	406376		ng/ul	99
89) Di-n-octyl phthalate	23.012	149	492531		ng/ul	100
90) Benzo(b)fluoranthene91) Benzo(k)fluoranthene	24.205	252	441637	31.09	•	99
93) Benzo(a)pyrene	24.275	252	398681	29.91		98
94) Indeno(1,2,3-cd)pyrene	25.133 29.193	252	413209	30.54		99
95) Dibenzo(a,h)anthracene	29.193	276	464926 390675	30.87		97 07
96) Benzo(g,h,i)perylene	30.421	278 276	385200m >	30.66		97 11/11/2/74
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^(#) = qualifier out of range (m) = manual integration (+) = signals summed