

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082718\
 Data File : BG036425.D
 Acq On : 27 Aug 2018 10:53
 Operator : JU/SJ
 Sample : PB112371BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB112371BS

Manual Integrations
 APPROVED

Sohil
 8/28/2018 12:12:42 PM

Quant Time: Aug 28 11:34:13 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 23 12:07:02 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	60243	20.00	ng	0.00
21) Naphthalene-d8	10.87	136	288433	20.00	ng	0.00
38) Acenaphthene-d10	14.69	164	189476	20.00	ng	0.00
63) Phenanthrene-d10	17.44	188	471562	20.00	ng	0.00
75) Chrysene-d12	21.70	240	433145	20.00	ng	0.00
86) Perylene-d12	24.92	264	454008	20.00	ng	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.63	112	418262	110.14	ng	0.00
7) Phenol-d6	7.22	99	577526	106.21	ng	0.00
23) Nitrobenzene-d5	9.23	82	326889	61.49	ng	0.00
41) 2,4,6-Tribromophenol	16.18	330	245165	108.44	ng	0.00
44) 2-Fluorobiphenyl	13.32	172	866385	65.29	ng	0.00
78) Terphenyl-d14	20.04	244	1318695	71.16	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.54	88	47993	27.079	ng	97
3) Pyridine	3.93	79	149859	30.123	ng	92
4) n-Nitrosodimethylamine	3.84	42	77486	34.969	ng	99
6) Aniline	7.38	93	228246	33.071	ng	98
8) 2-Chlorophenol	7.63	128	168911	41.014	ng	96
9) Benzaldehyde	7.20	77	110310	29.461	ng	95
10) Phenol	7.25	94	238028	41.832	ng	99
11) bis(2-Chloroethyl)ether	7.48	93	163440	36.231	ng	96
12) 1,3-Dichlorobenzene	7.95	146	165838	35.265	ng	97
13) 1,4-Dichlorobenzene	8.10	146	170395	35.889	ng	96
14) 1,2-Dichlorobenzene	8.42	146	162406	35.817	ng	97
15) Benzyl Alcohol	8.30	79	172374	39.325	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.60	45	307774	33.831	ng	96
17) 2-Methylphenol	8.50	107	161549	42.757	ng	99
18) Hexachloroethane	9.15	117	59013	34.667	ng	97
19) n-Nitroso-di-n-propylamine	8.87	70	145698	35.854	ng	96
20) 3+4-Methylphenols	8.83	107	221066	41.908	ng	99
22) Acetophenone	8.89	105	268270	35.419	ng	96
24) Nitrobenzene	9.27	77	199637	34.826	ng	96
25) Isophorone	9.79	82	414529	39.252	ng	97
26) 2-Nitrophenol	9.98	139	76348	33.610	ng	95
27) 2,4-Dimethylphenol	10.04	122	185508	44.110	ng	98
28) bis(2-Chloroethoxy)methane	10.28	93	239601	36.968	ng	98
29) 2,4-Dichlorophenol	10.52	162	189706	41.159	ng	98
30) 1,2,4-Trichlorobenzene	10.73	180	186755	34.636	ng	99
31) Naphthalene	10.93	128	540580	37.492	ng	98
32) Benzoic acid	10.17	122	110450m	36.597	ng	
33) 4-Chloroaniline	11.03	127	152781	23.124	ng	99
34) Hexachlorobutadiene	11.21	225	125136	34.542	ng	98
35) Caprolactam	11.80	113	73354m	39.951	ng	
36) 4-Chloro-3-methylphenol	12.14	107	213211	39.811	ng	98
37) 2-Methylnaphthalene	12.52	142	433986	41.136	ng	99
39) 1,2,4,5-Tetrachlorobenzene	12.89	216	244725	36.497	ng	99
40) Hexachlorocyclopentadiene	12.87	237	288663	82.740	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082718\
 Data File : BG036425.D
 Acq On : 27 Aug 2018 10:53
 Operator : JU/SJ
 Sample : PB112371BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB112371BS

Manual Integrations
 APPROVED

Sohil
 8/28/2018 12:12:42 PM

Quant Time: Aug 28 11:34:13 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 23 12:07:02 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.12	196	164282	40.220	ng	98
43) 2,4,5-Trichlorophenol	13.19	196	184320	42.308	ng	98
45) 1,1'-Biphenyl	13.53	154	569249	36.193	ng	99
46) 2-Chloronaphthalene	13.57	162	438638	36.532	ng	99
47) 2-Nitroaniline	13.76	65	141135	37.432	ng	98
48) Acenaphthylene	14.41	152	752735	40.114	ng	100
49) Dimethylphthalate	14.15	163	609496	39.394	ng	99
50) 2,6-Dinitrotoluene	14.26	165	125554	39.437	ng	95
51) Acenaphthene	14.76	154	485605	40.407	ng	99
52) 3-Nitroaniline	14.59	138	98866	28.817	ng	96
53) 2,4-Dinitrophenol	14.79	184	94254	78.161	ng	97
54) Dibenzofuran	15.09	168	710424	37.732	ng	99
55) 4-Nitrophenol	14.89	139	210442	75.020	ng	96
56) 2,4-Dinitrotoluene	15.05	165	174458	42.045	ng	92
57) Fluorene	15.74	166	569120	40.434	ng	97
58) 2,3,4,6-Tetrachlorophenol	15.31	232	176434	43.104	ng	97
59) Diethylphthalate	15.51	149	597756	38.337	ng	99
60) 4-Chlorophenyl-phenylether	15.73	204	330091	37.979	ng	98
61) 4-Nitroaniline	15.75	138	135573	37.763	ng	93
62) Azobenzene	16.03	77	581485	36.653	ng	97
64) 4,6-Dinitro-2-methylphenol	15.81	198	75802	36.593	ng	93
65) n-Nitrosodiphenylamine	15.94	169	530919	37.891	ng	99
66) 4-Bromophenyl-phenylether	16.63	248	215225	38.983	ng	98
67) Hexachlorobenzene	16.74	284	213000	37.730	ng	99
68) Atrazine	16.89	200	208197	39.587	ng	99
69) Pentachlorophenol	17.08	266	252023	65.685	ng	97
70) Phenanthrene	17.48	178	945772	39.471	ng	99
71) Anthracene	17.57	178	952815	39.891	ng	98
72) Carbazole	17.84	167	819082	34.337	ng	99
73) Di-n-butylphthalate	18.39	149	934818	35.192	ng	100
74) Fluoranthene	19.48	202	1087892	37.239	ng	99
76) Benzidine	19.66	184	575240	41.626	ng	99
77) Pyrene	19.84	202	1057821	39.714	ng	99
79) Butylbenzylphthalate	20.73	149	399231	37.662	ng	92
80) Benzo(a)anthracene	21.68	228	1054090	40.170	ng	99
81) 3,3'-Dichlorobenzidine	21.60	252	253650	24.254	ng	99
82) Chrysene	21.75	228	981375	39.456	ng	99
83) Bis(2-ethylhexyl)phthalate	21.60	149	557003	37.550	ng	99
84) Di-n-octyl phthalate	22.82	149	934824	37.730	ng	97
85) Indeno(1,2,3-cd)pyrene	28.58	276	1203290	41.652	ng	99
87) Benzo(b)fluoranthene	23.90	252	1066299	42.295	ng	96
88) Benzo(k)fluoranthene	23.97	252	993409	40.333	ng	99
89) Benzo(a)pyrene	24.77	252	1015403	41.914	ng	99
90) Dibenzo(a,h)anthracene	28.65	278	972544	41.583	ng	96
91) Benzo(g,h,i)perylene	29.72	276	990470	42.143	ng	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082718\
Data File : BG036425.D
Acq On : 27 Aug 2018 10:53
Operator : JU/SJ
Sample : PB112371BS
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PB112371BS

Manual Integrations
APPROVED

Sohil
8/28/2018 12:12:42 PM

Quant Time: Aug 28 11:34:13 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 23 12:07:02 2018
Response via : Initial Calibration

