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Data Path : Z:\svoasrv\HPCHEM1	\BNA_G\Da	ata\BG:	10321\		
Data File : BG050857.D		Instrument : BNA G			
Acq On : 3 Nov 2021 19:40					ClientSampleId :
Operator : CG/JU					SLCS426
Sample : PB140426BS					
Misc : ALS Vial : 42 Sample Multip]	1 d a m a 1				Manual IntegrationsAPPROVED
ALS Vial : 42 Sample Multip	lier: i				
Quant Time: Nov 07 07:42:35 202	01				Reviewed By :Jagrut Upadhyay 11/08/2021
Quant Method : Z:\svoasrv\HPCHE		i\Methc	ds\SFAM-EF	PA-BG110321.M	Reviewed By :Jagrut Upadhyay 11/08/2021 Supervised By :mohammad ahmed 11/08/2021
Quant Title : SVOA CALIBRATION	1	•			
QLast Update : Tue Nov 02 14:49					
Response via : Initial Calibrat	ion				
Common L			_		
Compound				Conc Units Dev	
Internal Standards					
	8.227	152	30092	20.000 ng/ul	0.00
20) Naphthalene-d8	11.053		136290	20.000 ng/ul	
38) Acenaphthene-d10	14.854		95346	20.000 ng/ul	0.00
64) Phenanthrene-d10	17.598		214591	20.000 ng/ul	0.00
79) Chrysene-d12	21.893	240	181704	20.000 ng/ul	
88) Perylene-d12	25.289	264	180428	20.000 ng/ul	
System Monitoring Compounds	_				
3) 1,4-Dioxane-d8	3.591		5330	5.717 ng/uL	0.00
4) Pyridine-d5	4.008		80908	29.007 ng/ul	0.00
7) Phenol-d5	7.375		98363	30.640 ng/ul	0.00
<pre>9) Bis-(2-Chloroethyl)eth 11) 2-Chlorophenol-d4</pre>	7.545 7.757		63610	30.674 ng/ul	0.00
15) 4-Methylphenol-d8	8.926		69390 77618	31.189 ng/ul 30.712 ng/ul	0.00 0.00
21) Nitrobenzene-d5	9.402		37376	32.271 ng/ul	0.00
24) 2-Nitrophenol-d4	10.125		41532	32.250 ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.671		70895	32.680 ng/ul	0.00
31) 4-Chloroaniline-d4	11.188		101060	30.762 ng/ul	0.00
46) Dimethylphthalate-d6	14.249	166	242996	33.312 ng/ul	0.00
49) Acenaphthylene-d8	14.549	160	295510	32.516 ng/ul	0.00
54) 4-Nitrophenol-d4	15.036		43187	32.653 ng/ul	0.00
60) Fluorene-d10	15.841		211636	32.751 ng/ul	0.00
65) 4,6-Dinitro-2-methylph			41820	32.144 ng/ul	0.00
73) Anthracene-d10 81) Pyrene-d10	17.698 19.972		321129	31.652 ng/ul	0.00
92) Benzo(a)pyrene-d12	25.060		382100 318658	32.558 ng/ul 31.948 ng/ul	0.00
	23.000	204	510050	51.540 hg/u1	-0.01
Target Compounds				Ova	alue
2) 1,4-Dioxane	3.626	88	12693	12.395 ng/uL	97
5) Pyridine	4.032	79	91018	31.524 ng/ul	97
6) Benzaldehyde	7.363	77	73760	36.428 ng/ul	96
8) Phenol	7.398	94	109620	33.009 ng/ul	99
10) Bis(2-Chloroethyl)ether	7.639	93	84009	33.797 ng/ul	99
12) 2-Chlorophenol 13) 2-Methylphenol	7.786	128	75428	33.391 ng/ul	98
14) 2,2'-oxybis(1-Chloropr	8.662 8.750	108 45	80952 128849	32.980 ng/ul 32.909 ng/ul	95
16) Acetophenone	9.055	105	129899	33.086 ng/ul	99 96
17) N-Nitroso-di-n-propyla	9.032	70	78809	33.270 ng/ul	99
18) 4-Methylphenol	8.991	108	87715	33.562 ng/ul	97
19) Hexachloroethane	9.320	117	31248	33.086 ng/ul	94
22) Nitrobenzene	9.443	77	112377	34.789 ng/ul	98
23) Isophorone	9.966	82	222286	35.457 ng/ul	98
25) 2-Nitrophenol	10.160	139	46154	35.723 ng/ul	97
26) 2,4-Dimethylphenol	10.207	107	96282	33.861 ng/ul	97
<pre>27) Bis(2-Chloroethoxy)met 29) 2,4-Dichlorophenol</pre>	10.442	93 160	119673	35.429 ng/ul	97
30) Naphthalene	10.694 11.106	162 128	76071 256136	35.950 ng/ul 34.368 ng/ul	95 98
32) 4-Chloroaniline	11.212	128	109247	33.490 ng/ul	98 99
33) Hexachlorobutadiene	11.376	225	47282	34.044 ng/ul	98
34) Caprolactam	11.969	113	31784	35.381 ng/ul	97
35) 4-Chloro-3-methylphenol	12.310	107	99083	36.655 ng/ul	98

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Data Path : Z:\	svoasrv\HPCHEM1	BNA G\Da	ata\BG:	110321\		
Data File : BG0						Instrument :
Acq On : 3	Nov 2021 19:40					BNA_G
Operator : CG/	JU					ClientSampleId :
	40426BS					SLCS426
Misc :		_				Manual Intervetion & DDD OV/CD
ALS Vial : 42	Sample Multipl	ier: 1				Manual IntegrationsAPPROVED
Quant Time: Nov	07 07:42:35 202	1				
	Z:\svoasrv\HPCHE		Mothe	de SEAM ED	A PC110201 M	Reviewed By :Jagrut Upadhyay 11/08/2021
Quant Title :	SVOA CALIBRATION	.mr./pinw_G	Ane cho	JUS (SFAM-EP	A-DG110521.M	Supervised By :mohammad ahmed 11/08/2021
	Tue Nov 02 14:49					
	Initial Calibrat					
Compoun	d	рт	OTen	Posponso	Conc Units Dev	////:->
	u 					·····
36) 2-Methyln	aphthalene	12.698	142	178483	35.154 ng/ul	100
37) 1-Methyln		12.910	142	177792	34.559 ng/ul	96
	etrachloroben			95100	34.231 ng/ul	96
	ocyclopentadiene	13.027	237	38145	28.592 ng/ul	98
41) 2,4,6-Tri		13.291		65340	35.943 ng/ul	96
42) 2,4,5-Tri		13.368		68809	35.256 ng/ul	96
43) 1,1'-Biph	•	13.691		240710	34.539 ng/ul	99
44) 2-Chloron 45) 2-Nitroan		13.738		188548	34.524 ng/ul	98
47) Dimethylp		13.938 14.296		76792 259731	35.391 ng/ul 35.610 ng/ul	94
48) 2,6-Dinit		14.425		56083	36.738 ng/ul	98 96
50) Acenaphthy		14.578		318147	34.929 ng/ul	99
51) 3-Nitroan		14.754		57086	36.156 ng/ul	91
52) Acenaphthe	ene	14.919	153	209942	35.053 ng/ul	95
53) 2,4-Dinitr	rophenol	14.966	184	27671	32.858 ng/ul	95
55) 4-Nitrophe		15.054	109	43780	36.092 ng/ul	95
56) Dibenzofur		15.248	168	300257	35.024 ng/ul	98
57) 2,4-Dinitr		15.207		79612	36.542 ng/ul#	
	etrachlorophenol		232	57240	37.320 ng/ul	99
59) Diethylpht 61) Fluorene	Indiate	15.648	149	279475	35.797 ng/ul	99
	enyl-phenyle	15.900 15.883	166 204	239503 123747	35.297 ng/ul	99 97
63) 4-Nitroani		15.918	138	58534	35.023 ng/ul 37.370 ng/ul	97
	o-2-methylph		198	46040	36.288 ng/ul#	
67) N-Nitrosod		16.094		214593	35.776 ng/ul	99
68) 4-Bromophe	nyl-phenylether	16.776	248	76704	35.934 ng/ul	98
69) Hexachloro	benzene	16.899	284	78286	35.674 ng/ul	99
70) Atrazine		17.034	200	89239	35.083 ng/ul	99
71) Pentachlor		17.246	266	35214	34.952 ng/ul	97
72) Phenanthre		17.639		407639	35.587 ng/ul	99
74) Anthracene		17.733	178	395762	34.434 ng/ul	99
76) Pentachlor	trachloroben	13.656 15.166	216 250	99876 91568	34.182 ng/uL 33.822 ng/uL	97 99
77) Carbazole	obenzene	17.998	167	380248	36.912 ng/ul	98
78) Di-n-butyl	phthalate	18.532		482494	35.645 ng/ul	100
80) Fluoranthe	•	19.637	202	493124	35.012 ng/ul	98
82) Pyrene		20.001	202	478338	34.758 ng/ul	99
83) Butylbenzy	•	20.865	149	211206	35.689 ng/ul	99
84) 3,3'-Dichl		21.782	252	148050	33.457 ng/ul	99
85) Benzo(a)an		21.875	228	445687	35.431 ng/ul	99
	lhexyl)phtha	21.746	149	302477	35.607 ng/ul	97
87) Chrysene	nhthalate	21.946	228 140	422253	35.139 ng/ul	99
89) Di-n-octyl 90) Benzo(b)fl		23.015 24.208	149 252	517018 448852	35.198 ng/ul	100
91) Benzo(k)fl		24.208	252 252	448852 409302	34.920 ng/ul 33.935 ng/ul	100 98
93) Benzo(a)py		25.136	252	419559	34.272 ng/ul	98
94) Indeno(1,2		29.196	276	475610	34.900 ng/ul	97
95) Dibenzo(a,		29.267		397487	34.472 ng/ul	98
96) Benzo(g,h,	i)perylene	30.424	276	391931m 🍗	34.358 ng/ul>	11/13/21 JU

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