

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
 Data File : BG025664.D
 Acq On : 26 Jan 2017 13:11
 Operator : SJ/MA
 Sample : I1387-08DL 20X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DA9R4DL

Manual Integrations
 APPROVED

mohammad
 1/27/2017 3:05:22 PM

Quant Time: Jan 27 10:10:29 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jan 27 05:43:39 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.18	152	67561	20.00	ng/ul	0.00
18) Naphthalene-d8	11.00	136	306382	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.81	164	256114	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.56	188	547540	20.00	ng/ul	0.00
75) Chrysene-d12	21.85	240	663498	20.00	ng/ul	0.00
83) Perylene-d12	25.24	264	694392	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.52	96	303	0.22	ng/uL	0.00
5) Phenol-d5	7.40	99	363	0.06	ng/ul	0.06
7) Bis-(2-Chloroethyl)ether-d	7.51	67	7648	2.05	ng/ul	0.02
9) 2-Chlorophenol-d4	7.74	132	4750	1.16	ng/ul	0.03
13) 4-Methylphenol-d8	8.95	113	2754m	0.56	ng/ul	0.06
19) Nitrobenzene-d5	9.39	128	2113	0.97	ng/ul	0.04
22) 2-Nitrophenol-d4	10.11	143	4573	1.76	ng/ul	0.03
26) 2,4-Dichlorophenol-d3	10.69	165	6888m	1.35	ng/ul	0.06
29) 4-Chloroaniline-d4	11.16	131	92	0.02	ng/ul	0.00
43) Dimethylphthalate-d6	14.21	166	36315	2.01	ng/ul	0.00
46) Acenaphthylene-d8	14.51	160	45792	2.34	ng/ul	0.00
51) 4-Nitrophenol-d4	15.04	143	250	0.10	ng/ul	0.00
57) Fluorene-d10	15.80	176	37533	2.20	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul	
70) Anthracene-d10	17.66	188	56566m	2.43	ng/ul	0.00
76) Pyrene-d10	19.94	212	68525	2.48	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.01	264	64043	2.19	ng/ul	0.02

Target Compounds						Qvalue
55) 2,3,4,6-Tetrachlorophenol	15.45	232	14857	2.87	ng/ul#	87
68) Pentachlorophenol	17.22	266	258122	85.93	ng/ul	91

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

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