

Data Path : U:\HPCHEM1\BNA G\DATA\BG082117\
 Data File : BG028400.D
 Acq On : 21 Aug 2017 22:40
 Operator : SJ/JU
 Sample : I4827-16
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 B0AF4

Quant Time: Aug 22 13:57:59 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG080217.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 22 04:33:26 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.35	152	38966	20.00	ng/ul	0.00
18) Naphthalene-d8	11.16	136	167539	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.95	164	117221	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.70	188	334570	20.00	ng/ul	0.00
75) Chrysene-d12	22.02	240	382457	20.00	ng/ul	0.00
83) Perylene-d12	25.53	264	378052	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.83	96	4294	5.13	ng/uL	0.00
5) Phenol-d5	7.49	99	101225	23.65	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.65	67	66205	24.43	ng/ul	0.00
9) 2-Chlorophenol-d4	7.88	132	72513	26.86	ng/ul	0.00
13) 4-Methylphenol-d8	9.04	113	81923	22.90	ng/ul	0.00
19) Nitrobenzene-d5	9.51	128	37370	28.53	ng/ul	0.00
22) 2-Nitrophenol-d4	10.24	143	45507	31.68	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.78	165	83532	28.11	ng/ul	0.00
29) 4-Chloroaniline-d4	11.30	131	86856	26.26	ng/ul	0.00
43) Dimethylphthalate-d6	14.34	166	291611	27.97	ng/ul	0.00
46) Acenaphthylene-d8	14.65	160	340686	28.98	ng/ul	0.00
51) 4-Nitrophenol-d4	15.15	143	44247	26.69	ng/ul	0.00
57) Fluorene-d10	15.94	176	265587	28.39	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	16.06	200	53306	24.53	ng/ul	0.00
70) Anthracene-d10	17.79	188	482998	30.11	ng/ul	0.00
76) Pyrene-d10	20.07	212	583352	29.74	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.29	264	519262	29.34	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	7.52	94	6657	1.565	ng/ul 93
44) Dimethylphthalate	14.39	163	66816	6.960	ng/ul 100
81) Bis(2-ethylhexyl)phthalate	21.86	149	13613	1.120	ng/ul 96

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

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