

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100818\
 Data File : BM017072.D
 Acq On : 08 Oct 2018 18:51
 Operator : MJ/SJ
 Sample : J5234-18
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AK3

Quant Time: Oct 09 04:02:16 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 09 04:00:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	196712	20.00	ng/ul	0.00
18) Naphthalene-d8	10.47	136	859508	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.33	164	466182	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.08	188	1044726	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	1165968	20.00	ng/ul	0.00
86) Perylene-d12	23.49	264	1258683	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.25	96	20668	5.22	ng/uL	0.00
5) Phenol-d5	6.87	99	108028	6.57	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.03	67	275969	27.68	ng/ul	0.00
9) 2-Chlorophenol-d4	7.23	132	314815	23.03	ng/ul	0.00
13) 4-Methylphenol-d8	8.39	113	194114	15.07	ng/ul	0.00
19) Nitrobenzene-d5	8.85	128	182401	29.04	ng/ul	0.00
22) 2-Nitrophenol-d4	9.56	143	170179	26.63	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.09	165	342768	26.07	ng/ul	0.00
29) 4-Chloroaniline-d4	10.62	131	21502	1.36	ng/ul	0.00
44) Dimethylphthalate-d6	13.75	166	1102977	29.82	ng/ul	0.00
47) Acenaphthylene-d8	14.02	160	1448322	30.63	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	16992	2.43	ng/ul	0.00
58) Fluorene-d10	15.33	176	1007992	30.60	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.45	200	92290	14.98	ng/ul	0.00
71) Anthracene-d10	17.18	188	1535489	30.77	ng/ul	0.00
79) Pyrene-d10	19.47	212	1700500	32.33	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.35	264	1980612	30.74	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.89	94	36789	2.236 ng/ul	99
45) Dimethylphthalate	13.79	163	79191	2.162 ng/ul	98

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

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