

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060818\
 Data File : BM015566.D
 Acq On : 08 Jun 2018 16:35
 Operator : SJ/JU
 Sample : J3261-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C07Q3

Quant Time: Jun 08 17:50:31 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM060818MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 08 14:39:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.35	152	73848	20.00	ng/ul	0.00
18) Naphthalene-d8	11.18	136	372032	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.96	164	240070	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.69	188	599679	20.00	ng/ul	0.00
77) Chrysene-d12	21.79	240	717007	20.00	ng/ul	0.00
85) Perylene-d12	24.20	264	814292	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.55	96	9244	6.08	ng/uL	0.00
5) Phenol-d5	7.49	99	45077	6.72	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.66	67	153112	37.80	ng/ul	0.00
9) 2-Chlorophenol-d4	7.87	132	151428	29.54	ng/ul	0.00
13) 4-Methylphenol-d8	9.06	113	97250	16.84	ng/ul	0.00
19) Nitrobenzene-d5	9.53	128	99515	39.78	ng/ul	0.00
22) 2-Nitrophenol-d4	10.26	143	110026	40.89	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.79	165	186456	33.47	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
43) Dimethylphthalate-d6	14.36	166	873582	43.95	ng/ul	0.00
46) Acenaphthylene-d8	14.66	160	930735	43.42	ng/ul	0.00
51) 4-Nitrophenol-d4	15.16	143	26708	6.96	ng/ul	0.00
57) Fluorene-d10	15.94	176	709465	42.18	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	16.07	200	129401	35.82	ng/ul	0.00
70) Anthracene-d10	17.78	188	1187132	43.42	ng/ul	0.00
78) Pyrene-d10	20.03	212	1400714	45.42	ng/ul	0.00
89) Benzo(a)pyrene-d12	24.05	264	1785726	43.75	ng/ul	0.00

Target Compounds				Ovalue		
10) 2-Chlorophenol	7.90	128	40303	7.630	ng/ul#	89
20) Nitrobenzene	9.57	77	878885	130.072	ng/ul	91
23) 2-Nitrophenol	10.29	139	9486	3.058	ng/ul#	87
36) 1,2,4,5-Tetrachlorobenzene	13.17	216	12265	1.770	ng/ul	97
68) Pentachlorophenol	17.34	266	4246	1.036	ng/ul#	90
72) 1,2,3,4-Tetrachlorobenzene	13.77	216	99812	13.801	ng/uL	98

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

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