

Data Path : Z:\HPCHEM1\BNA M\DATA\BM100517\
 Data File : BM011927.D
 Acq On : 05 Oct 2017 15:55
 Operator : SJ/JU
 Sample : I5572-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Oct 06 04:01:59 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM100317.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 06 04:00:01 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.87	152	251527	20.00	ng/ul	0.00
18) Naphthalene-d8	10.67	136	1021434	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.52	164	593567	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.26	188	1246779	20.00	ng/ul	0.00
75) Chrysene-d12	21.43	240	1081592	20.00	ng/ul	0.00
83) Perylene-d12	23.76	264	1165614	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.29	96	29664	5.58	ng/uL	0.00
5) Phenol-d5	7.03	99	27815	1.29	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.20	67	78779	6.15	ng/ul	0.00
9) 2-Chlorophenol-d4	7.40	132	91140	5.25	ng/ul	0.00
13) 4-Methylphenol-d8	8.57	113	56683	3.03	ng/ul	0.00
19) Nitrobenzene-d5	9.03	128	49785	6.84	ng/ul	0.00
22) 2-Nitrophenol-d4	9.76	143	55669	6.79	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.30	165	97297	5.88	ng/ul	0.00
29) 4-Chloroaniline-d4	10.83	131	8394	0.46	ng/ul	0.02
43) Dimethylphthalate-d6	13.92	166	393815	7.66	ng/ul	0.00
46) Acenaphthylene-d8	14.21	160	426250	7.05	ng/ul	0.00
51) 4-Nitrophenol-d4	14.75	143	1328	0.15	ng/ul	0.04
57) Fluorene-d10	15.50	176	310133	6.83	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.63	200	36205	4.32	ng/ul	0.00
70) Anthracene-d10	17.36	188	490158	7.98	ng/ul	0.00
76) Pyrene-d10	19.64	212	494970	10.15	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.60	264	426709	7.61	ng/ul	0.00

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

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