

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070916\
 Data File : BM006389.D
 Acq On : 09 Jul 2016 22:05
 Operator : UM/SJ
 Sample : H3914-13
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Jul 11 04:58:36 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM070216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 11 04:57:47 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	144398	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	612171	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.28	164	383469	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.03	188	918166	20.00	ng/ul	0.00
75) Chrysene-d12	21.23	240	1097150	20.00	ng/ul	0.00
83) Perylene-d12	23.44	264	1154952	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.15	96	1638	0.48	ng/uL	0.00
5) Phenol-d5	6.81	99	83023	6.14	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.97	67	209645	26.28	ng/ul	0.00
9) 2-Chlorophenol-d4	7.16	132	225254	21.77	ng/ul	0.00
13) 4-Methylphenol-d8	8.34	113	151180	13.82	ng/ul	0.00
19) Nitrobenzene-d5	8.79	128	137123	28.11	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	155665	28.15	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.04	165	254851	25.44	ng/ul	0.00
29) 4-Chloroaniline-d4	10.56	131	11778	1.13	ng/ul	0.00
43) Dimethylphthalate-d6	13.69	166	1003382	31.35	ng/ul	0.00
46) Acenaphthylene-d8	13.97	160	1163172	29.36	ng/ul	0.00
51) 4-Nitrophenol-d4	14.50	143	47965	7.91	ng/ul	0.00
57) Fluorene-d10	15.28	176	882102	31.97	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.41	200	191657	35.63	ng/ul	0.00
70) Anthracene-d10	17.13	188	1576356	35.98	ng/ul	0.00
76) Pyrene-d10	19.43	212	1817695	34.74	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.30	264	1990599	37.10	ng/ul	0.00

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

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