

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM072716\
 Data File : BM006705.D
 Acq On : 28 Jul 2016 01:57
 Operator : UM/SJ
 Sample : H4192-09
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA0E4

Quant Time: Jul 28 11:01:30 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 28 01:25:43 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	159438	20.00	ng/ul	0.00
18) Naphthalene-d8	10.39	136	685776	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.26	164	400557	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.01	188	917346	20.00	ng/ul	0.00
75) Chrysene-d12	21.22	240	1069042	20.00	ng/ul	0.00
83) Perylene-d12	23.42	264	1153475	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.13	96	2879	0.70	ng/uL	0.00
5) Phenol-d5	6.79	99	100751	7.19	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.95	67	269894	31.36	ng/ul	0.00
9) 2-Chlorophenol-d4	7.14	132	276119	25.12	ng/ul	0.00
13) 4-Methylphenol-d8	8.32	113	180282	16.40	ng/ul	0.00
19) Nitrobenzene-d5	8.76	128	166188	31.39	ng/ul	0.00
22) 2-Nitrophenol-d4	9.47	143	184860	32.32	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.02	165	297053	28.33	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
43) Dimethylphthalate-d6	13.67	166	1028163	32.17	ng/ul	0.00
46) Acenaphthylene-d8	13.95	160	1276687	31.77	ng/ul	0.00
51) 4-Nitrophenol-d4	14.50	143	28368	5.16	ng/ul	0.00
57) Fluorene-d10	15.26	176	905686	32.67	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.40	200	136187	27.24	ng/ul	0.00
70) Anthracene-d10	17.11	188	1396248	32.62	ng/ul	0.00
76) Pyrene-d10	19.42	212	1595890	32.18	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.28	264	1657474	31.75	ng/ul	0.00

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

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