

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM073016\
 Data File : BM006759.D
 Acq On : 30 Jul 2016 05:45
 Operator : UM/SJ
 Sample : H4192-09
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA0E4

Quant Time: Aug 01 04:23:48 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 30 01:37:26 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	137662	20.00	ng/ul	0.00
7) Naphthalene-d8	10.38	136	583468	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.26	164	346545	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.01	188	795413	20.00	ng/ul	0.00
29) Chrysene-d12	21.21	240	952619	20.00	ng/ul	0.00
34) Perylene-d12	23.41	264	1014174	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.13	96	2944	0.86	ng/uL	0.00
3) Phenol-d5	6.79	99	91035	7.77	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.95	67	245600	33.85	ng/ul	0.00
5) 2-Chlorophenol-d4	7.14	132	250991	26.82	ng/ul	0.00
6) 4-Methylphenol-d8	8.32	113	162501	17.89	ng/ul	0.00
8) Nitrobenzene-d5	8.76	128	149699	33.89	ng/ul	0.00
9) 2-Nitrophenol-d4	9.47	143	166049	35.35	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.01	165	265952	29.97	ng/ul	0.00
12) 4-Chloroaniline-d4	10.39	131	293	0.03	ng/ul	-0.14
16) Dimethylphthalate-d6	13.67	166	948548	34.91	ng/ul	0.00
17) Acenaphthylene-d8	13.94	160	1166632	33.64	ng/ul	0.00
20) 4-Nitrophenol-d4	14.49	143	25785	5.46	ng/ul	0.00
21) Fluorene-d10	15.26	176	830971	34.20	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.40	200	131336	31.81	ng/ul	0.00
26) Anthracene-d10	17.11	188	1290083	34.16	ng/ul	0.00
30) Pyrene-d10	19.41	212	1485989	34.23	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.27	264	1613135	34.58	ng/ul	0.00

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

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