

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM073016\
 Data File : BM006740.D
 Acq On : 29 Jul 2016 17:23
 Operator : UM/SJ
 Sample : PB92420BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK20

Quant Time: Jul 30 01:38:06 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	98677	20.00	ng/ul	0.00
7) Naphthalene-d8	10.38	136	425907	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.26	164	248390	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.01	188	578674	20.00	ng/ul	0.00
29) Chrysene-d12	21.21	240	701561	20.00	ng/ul	0.00
34) Perylene-d12	23.41	264	737366	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.13	96	12092	4.96	ng/uL	0.00
3) Phenol-d5	6.79	99	287540	34.25	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.95	67	188313	36.21	ng/ul	0.00
5) 2-Chlorophenol-d4	7.14	132	226416	33.75	ng/ul	0.00
6) 4-Methylphenol-d8	8.32	113	225264	34.60	ng/ul	0.00
8) Nitrobenzene-d5	8.76	128	109370	33.92	ng/ul	0.00
9) 2-Nitrophenol-d4	9.47	143	116736	34.05	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.02	165	201704	31.13	ng/ul	0.00
12) 4-Chloroaniline-d4	10.53	131	301199	42.21	ng/ul	0.00
16) Dimethylphthalate-d6	13.67	166	739599	37.98	ng/ul	0.00
17) Acenaphthylene-d8	13.95	160	871943	35.08	ng/ul	0.00
20) 4-Nitrophenol-d4	14.49	143	111355	32.92	ng/ul	0.00
21) Fluorene-d10	15.26	176	623795	35.81	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.40	200	96232	32.04	ng/ul	0.00
26) Anthracene-d10	17.11	188	1035963	37.70	ng/ul	0.00
30) Pyrene-d10	19.42	212	1197738	37.47	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.27	264	1311943	38.68	ng/ul	0.00

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

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