

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060616\
 Data File : BM005672.D
 Acq On : 06 Jun 2016 16:30
 Operator : UM/SJ
 Sample : PB91122BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK22

Quant Time: Jun 07 04:26:38 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM060216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 07 04:25:20 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	96398	20.00	ng/ul	0.00
7) Naphthalene-d8	10.55	136	421379	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.40	164	203660	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.14	188	387318	20.00	ng/ul	0.00
29) Chrysene-d12	21.33	240	361930	20.00	ng/ul	0.00
34) Perylene-d12	23.59	264	370501	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.22	96	7547	3.29	ng/uL	0.00
3) Phenol-d5	6.93	99	280820	33.14	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.09	67	179798	34.86	ng/ul	0.00
5) 2-Chlorophenol-d4	7.29	132	234594	33.55	ng/ul	0.00
6) 4-Methylphenol-d8	8.47	113	220875	32.95	ng/ul	0.00
8) Nitrobenzene-d5	8.92	128	110519	34.16	ng/ul	0.00
9) 2-Nitrophenol-d4	9.64	143	117222	33.86	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.18	165	196942	32.01	ng/ul	0.00
12) 4-Chloroaniline-d4	10.69	131	305705	41.81	ng/ul	0.00
16) Dimethylphthalate-d6	13.81	166	562821	35.95	ng/ul	0.00
17) Acenaphthylene-d8	14.09	160	716271	35.17	ng/ul	0.00
20) 4-Nitrophenol-d4	14.62	143	72116	32.17	ng/ul	0.00
21) Fluorene-d10	15.39	176	470755	35.01	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.53	200	45904	25.17	ng/ul	0.00
26) Anthracene-d10	17.24	188	669488	36.40	ng/ul	0.00
30) Pyrene-d10	19.54	212	667995	36.72	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.44	264	632682	37.11	ng/ul	0.00

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

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