

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040116\
 Data File : BM004859.D
 Acq On : 02 Apr 2016 02:41
 Operator : UM/SJ
 Sample : H2004-07
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AJ7

Quant Time: Apr 02 03:25:43 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM040116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Apr 02 00:53:13 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	52260	20.00	ng/ul	0.00
18) Naphthalene-d8	10.62	136	245166	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.46	164	153661	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.20	188	372695	20.00	ng/ul	0.00
78) Chrysene-d12	21.39	240	456026	20.00	ng/ul	-0.01
86) Perylene-d12	23.68	264	474055	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.32	96	1476	1.14	ng/uL	0.00
5) Phenol-d5	6.99	99	30865	6.65	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.16	67	90712	36.05	ng/ul	0.00
9) 2-Chlorophenol-d4	7.36	132	97170	27.57	ng/ul	0.00
13) 4-Methylphenol-d8	8.52	113	62502	16.03	ng/ul	0.00
19) Nitrobenzene-d5	8.98	128	62454	35.90	ng/ul	0.00
22) 2-Nitrophenol-d4	9.70	143	70453	36.43	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.23	165	113430	31.46	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
44) Dimethylphthalate-d6	13.87	166	459674	37.92	ng/ul	0.00
47) Acenaphthylene-d8	14.15	160	540973	37.30	ng/ul	0.00
52) 4-Nitrophenol-d4	14.66	143	12212	5.10	ng/ul	0.00
58) Fluorene-d10	15.45	176	399997	38.18	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.57	200	72451	32.44	ng/ul	0.00
71) Anthracene-d10	17.30	188	636601	38.17	ng/ul	0.00
79) Pyrene-d10	19.60	212	741782	37.47	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.54	264	878311	41.82	ng/ul	0.00

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

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