

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012316\
 Data File : BM004120.D
 Acq On : 22 Jan 2016 21:13
 Operator : SJ/UM
 Sample : H1174-15
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BC9K1

Quant Time: Jan 23 02:40:44 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM012016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jan 23 02:35:35 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	19522	20.00	ng/ul	0.00
18) Naphthalene-d8	10.45	136	95054	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.32	164	61990	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	155459	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	212964	20.00	ng/ul	0.00
86) Perylene-d12	23.51	264	239495	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.12	96	457	1.24	ng/uL	0.00
5) Phenol-d5	6.84	99	12815	6.73	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.99	67	29155	28.76	ng/ul	0.00
9) 2-Chlorophenol-d4	7.19	132	35538	23.97	ng/ul	0.00
13) 4-Methylphenol-d8	8.37	113	24420	14.95	ng/ul	0.00
19) Nitrobenzene-d5	8.82	128	20355	28.89	ng/ul	0.00
22) 2-Nitrophenol-d4	9.54	143	21967	28.10	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.07	165	37198	25.60	ng/ul	0.00
29) 4-Chloroaniline-d4	10.61	131	243	0.13	ng/ul	0.02
44) Dimethylphthalate-d6	13.73	166	172328	32.31	ng/ul	0.00
47) Acenaphthylene-d8	14.01	160	196146	31.59	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	4622	4.33	ng/ul	0.01
58) Fluorene-d10	15.32	176	158200	34.58	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.45	200	21887	22.84	ng/ul	0.00
71) Anthracene-d10	17.17	188	285560	37.76	ng/ul	0.00
79) Pyrene-d10	19.47	212	350821	36.77	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.37	264	461040	36.91	ng/ul	0.00

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

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