

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM120717\
 Data File : BM013259.D
 Acq On : 08 Dec 2017 00:38
 Operator : SJ/JU
 Sample : I6719-14
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JHPH0

Quant Time: Dec 08 07:05:31 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 08 03:46:15 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	191382	20.00	ng/ul	0.00
18) Naphthalene-d8	10.47	136	729999	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.34	164	388671	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.09	188	836252	20.00	ng/ul	0.00
75) Chrysene-d12	21.27	240	863820	20.00	ng/ul	0.00
83) Perylene-d12	23.50	264	871118	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.24	96	25234	6.53	ng/uL	0.00
5) Phenol-d5	6.88	99	542113	32.74	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.04	67	319213	33.80	ng/ul	0.00
9) 2-Chlorophenol-d4	7.23	132	460349	35.58	ng/ul	0.00
13) 4-Methylphenol-d8	8.40	113	443959	31.38	ng/ul	0.00
19) Nitrobenzene-d5	8.85	128	229953	39.72	ng/ul	0.00
22) 2-Nitrophenol-d4	9.57	143	251006	40.51	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.10	165	443485	34.73	ng/ul	0.00
29) 4-Chloroaniline-d4	10.62	131	560636	47.78	ng/ul	0.00
43) Dimethylphthalate-d6	13.75	166	1276169	36.90	ng/ul	0.00
46) Acenaphthylene-d8	14.03	160	1652489	38.72	ng/ul	0.00
51) 4-Nitrophenol-d4	14.55	143	204884	34.09	ng/ul	-0.01
57) Fluorene-d10	15.33	176	1083380	36.42	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.46	200	160395	30.49	ng/ul	0.00
70) Anthracene-d10	17.19	188	1619063	38.66	ng/ul	0.00
76) Pyrene-d10	19.49	212	1810103	42.06	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.37	264	1725244	38.88	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.90	94	44075	2.693 ng/ul	89
44) Dimethylphthalate	13.80	163	315149	9.527 ng/ul	99

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

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