

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111119\
 Data File : BM023647.D
 Acq On : 11 Nov 2019 21:29
 Operator : JU
 Sample : K5559-02
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0CP6

Quant Time: Nov 12 01:16:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111119MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 12 00:02:36 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	338870	20.00	ng/ul	0.05
18) Naphthalene-d8	10.35	136	1784382	20.00	ng/ul	0.03
36) Acenaphthene-d10	14.20	164	1193239	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.95	188	2618625	20.00	ng/ul	0.00
78) Chrysene-d12	21.15	240	1905118	20.00	ng/ul	0.00
86) Perylene-d12	23.31	264	1845829	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.08	96	54574	6.64	ng/uL	0.00
5) Phenol-d5	6.74	99	278016	9.00	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	6.91	67	773289	43.57	ng/ul	0.02
9) 2-Chlorophenol-d4	7.09	132	850592	38.01	ng/ul	0.01
13) 4-Methylphenol-d8	8.27	113	542271	22.01	ng/ul	0.00
19) Nitrobenzene-d5	8.71	128	557097	43.62	ng/ul	0.01
22) 2-Nitrophenol-d4	9.42	143	648451	48.30	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.96	165	1135318	37.09	ng/ul	0.01
29) 4-Chloroaniline-d4	10.47	131	87972	2.62	ng/ul	0.01
44) Dimethylphthalate-d6	13.63	166	4008889	43.08	ng/ul	0.00
47) Acenaphthylene-d8	13.89	160	4569696	43.03	ng/ul	0.00
52) 4-Nitrophenol-d4	14.42	143	129566	8.16	ng/ul	0.00
58) Fluorene-d10	15.20	176	3427766	41.71	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.33	200	646852	42.70	ng/ul	0.00
71) Anthracene-d10	17.05	188	5467935	44.02	ng/ul	0.00
79) Pyrene-d10	19.35	212	5589811	56.27	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.18	264	4485068	46.18	ng/ul	0.00

Target Compounds

					Ovalue	
20) Nitrobenzene	8.76	77	5970462	176.800	ng/ul	93
23) 2-Nitrophenol	9.45	139	51681	3.461	ng/ul	98
37) 1,2,4,5-Tetrachlorobenzene	12.38	216	1959960	50.270	ng/ul	99
45) Dimethylphthalate	13.67	163	228578	2.507	ng/ul	97
73) 1,2,3,4-Tetrachlorobenzene	13.00	216	6999465	188.541	ng/uL	99
74) Pentachlorobenzene	14.53	250	191267	4.663	ng/uL	99

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

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