

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120619\
 Data File : BM024016.D
 Acq On : 06 Dec 2019 15:07
 Operator : JU
 Sample : K6007-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0BL4

Quant Time: Dec 06 15:39:26 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 06 13:09:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	388425	20.00	ng/ul	0.00
18) Naphthalene-d8	10.25	136	1608803	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.14	164	966386	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.89	188	1878806	20.00	ng/ul	0.00
78) Chrysene-d12	21.10	240	1625309	20.00	ng/ul	0.00
86) Perylene-d12	23.24	264	1989587	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.03	96	32552	3.73	ng/uL	0.00
5) Phenol-d5	6.67	99	670642	19.70	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.83	67	397588	20.08	ng/ul	0.00
9) 2-Chlorophenol-d4	7.02	132	526918	20.02	ng/ul	0.00
13) 4-Methylphenol-d8	8.19	113	524260	19.52	ng/ul	0.00
19) Nitrobenzene-d5	8.63	128	247229	19.80	ng/ul	0.00
22) 2-Nitrophenol-d4	9.35	143	281910	19.86	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.88	165	521829	19.42	ng/ul	0.00
29) 4-Chloroaniline-d4	10.40	131	355852	12.06	ng/ul	0.00
44) Dimethylphthalate-d6	13.56	166	1636384	21.79	ng/ul	0.00
47) Acenaphthylene-d8	13.83	160	1972562	22.42	ng/ul	0.00
52) 4-Nitrophenol-d4	14.37	143	179967	14.26	ng/ul	0.00
58) Fluorene-d10	15.14	176	1396713	22.10	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.29	200	171409	14.53	ng/ul	0.00
71) Anthracene-d10	16.99	188	1940885	21.63	ng/ul	0.00
79) Pyrene-d10	19.30	212	1991575	23.71	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.11	264	2275971	21.64	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.69	94	53079	1.516	ng/ul 95
45) Dimethylphthalate	13.61	163	321258	4.371	ng/ul 99

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

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