

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120619\
 Data File : BM024018.D
 Acq On : 06 Dec 2019 16:18
 Operator : JU
 Sample : K6007-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0BL6

Quant Time: Dec 06 17:08:14 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	7.47	152	358506	20.00	ng/ul	0.00
18) Naphthalene-d8	10.25	136	1433987	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.14	164	842187	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.89	188	1755286	20.00	ng/ul	0.00
78) Chrysene-d12	21.10	240	1692454	20.00	ng/ul	0.00
86) Perylene-d12	23.24	264	2040863	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.03	96	39461	4.90	ng/uL	0.00
5) Phenol-d5	6.67	99	692497	22.04	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.83	67	441010	24.13	ng/ul	0.00
9) 2-Chlorophenol-d4	7.02	132	571200	23.52	ng/ul	0.00
13) 4-Methylphenol-d8	8.19	113	515611	20.80	ng/ul	0.00
19) Nitrobenzene-d5	8.63	128	266577	23.95	ng/ul	0.00
22) 2-Nitrophenol-d4	9.35	143	289799	22.91	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.88	165	522880	21.83	ng/ul	0.00
29) 4-Chloroaniline-d4	10.40	131	594240	22.60	ng/ul	0.00
44) Dimethylphthalate-d6	13.56	166	1661729	25.39	ng/ul	0.00
47) Acenaphthylene-d8	13.83	160	1981961	25.84	ng/ul	0.00
52) 4-Nitrophenol-d4	14.37	143	204686	18.60	ng/ul	0.00
58) Fluorene-d10	15.14	176	1438179	26.12	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.28	200	213547	19.38	ng/ul	0.00
71) Anthracene-d10	16.99	188	2203039	26.28	ng/ul	0.00
79) Pyrene-d10	19.30	212	2506138	28.66	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.11	264	3091240	28.66	ng/ul	0.00

Target Compounds					Ovalue	
6) Phenol	6.69	94	58011	1.795	ng/ul#	90
45) Dimethylphthalate	13.61	163	276088	4.310	ng/ul	99

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

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