

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120619\
 Data File : BM024014.D
 Acq On : 06 Dec 2019 13:55
 Operator : JU
 Sample : K6007-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0BL1

Quant Time: Dec 06 14:35:48 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	309941	20.00	ng/ul	0.00
18) Naphthalene-d8	10.25	136	1238351	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.14	164	741177	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.89	188	1525843	20.00	ng/ul	0.00
78) Chrysene-d12	21.10	240	1486520	20.00	ng/ul	0.00
86) Perylene-d12	23.24	264	1836327	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.03	96	29696	4.26	ng/uL	0.00
5) Phenol-d5	6.67	99	561818	20.68	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.83	67	346295	21.92	ng/ul	0.00
9) 2-Chlorophenol-d4	7.02	132	450792	21.47	ng/ul	0.00
13) 4-Methylphenol-d8	8.20	113	429497	20.04	ng/ul	0.00
19) Nitrobenzene-d5	8.63	128	207581	21.60	ng/ul	0.00
22) 2-Nitrophenol-d4	9.35	143	224656	20.56	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.89	165	408565	19.75	ng/ul	0.00
29) 4-Chloroaniline-d4	10.40	131	470042	20.70	ng/ul	0.00
44) Dimethylphthalate-d6	13.56	166	1267944	22.02	ng/ul	0.00
47) Acenaphthylene-d8	13.83	160	1516193	22.46	ng/ul	0.00
52) 4-Nitrophenol-d4	14.37	143	149679	15.46	ng/ul	0.00
58) Fluorene-d10	15.14	176	1134867	23.42	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.29	200	149534	15.61	ng/ul	0.00
71) Anthracene-d10	16.99	188	1702958	23.37	ng/ul	0.00
79) Pyrene-d10	19.30	212	1921809	25.02	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.11	264	2389030	24.61	ng/ul	0.00

Target Compounds				Ovalue		
6) Phenol	6.70	94	55394	1.982	ng/ul	93
45) Dimethylphthalate	13.61	163	418848	7.430	ng/ul	100
70) Phenanthrene	16.93	178	259966	3.016	ng/ul	98
77) Fluoranthene	18.96	202	493816	5.342	ng/ul	98
80) Pyrene	19.33	202	415184	4.145	ng/ul	97
83) Benzo(a)anthracene	21.09	228	227019	2.285	ng/ul	96
85) Chrysene	21.14	228	206551	2.214	ng/ul	97
88) Benzo(b)fluoranthene	22.61	252	298134	2.625	ng/ul	97
91) Benzo(a)pyrene	23.15	252	208822	1.945	ng/ul#	92
92) Indeno(1,2,3-cd)pyrene	25.36	276	141328	1.060	ng/ul	98
94) Benzo(g,h,i)perylene	26.01	276	137024	1.227	ng/ul	98

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120619\
Data File : BM024014.D
Acq On : 06 Dec 2019 13:55
Operator : JU
Sample : K6007-03
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0BL1

Quant Time: Dec 06 14:35:48 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

