

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012120\
 Data File : BM024561.D
 Acq On : 21 Jan 2020 13:56
 Operator : JU
 Sample : PB126224BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK24

Quant Time: Jan 21 15:28:38 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011520MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 21 11:19:17 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	216745	20.00	ng/ul	0.00
18) Naphthalene-d8	10.21	136	965299	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.10	164	691670	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.85	188	1581000	20.00	ng/ul	0.00
78) Chrysene-d12	21.06	240	1437562	20.00	ng/ul	0.00
86) Perylene-d12	23.17	264	1266129	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.06	96	29468	6.64	ng/uL	0.00
5) Phenol-d5	6.65	99	572514	35.63	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.80	67	327250	36.67	ng/ul	0.00
9) 2-Chlorophenol-d4	6.99	132	524513	38.23	ng/ul	0.00
13) 4-Methylphenol-d8	8.16	113	522376	35.26	ng/ul	0.00
19) Nitrobenzene-d5	8.59	128	268444	39.58	ng/ul	0.00
22) 2-Nitrophenol-d4	9.31	143	324382	44.16	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.84	165	586986	35.50	ng/ul	0.00
29) 4-Chloroaniline-d4	10.35	131	770601	41.92	ng/ul	0.00
44) Dimethylphthalate-d6	13.53	166	1963209	36.22	ng/ul	0.00
47) Acenaphthylene-d8	13.79	160	2352650	37.72	ng/ul	0.00
52) 4-Nitrophenol-d4	14.32	143	325143	35.95	ng/ul	0.00
58) Fluorene-d10	15.10	176	1724352	36.15	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.23	200	313725	34.76	ng/ul	0.00
71) Anthracene-d10	16.95	188	2703999	36.78	ng/ul	0.00
79) Pyrene-d10	19.26	212	3091324	40.23	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.04	264	2442049	37.19	ng/ul	0.00

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

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