

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120520\
 Data File : BM028048.D
 Acq On : 05 Dec 2020 15:22
 Operator : JU/CG
 Sample : PB133355BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK355

Quant Time: Dec 05 15:57:02 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120320.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Dec 05 11:13:40 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.29	152	203083	20.00	ng/ul	0.00
20) Naphthalene-d8	11.12	136	843186	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.91	164	489946	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.63	188	999812	20.00	ng/ul	0.00
79) Chrysene-d12	21.73	240	898144	20.00	ng/ul	0.00
88) Perylene-d12	24.14	264	849017	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.47	96	33565	6.17	ng/uL	0.00
4) Pyridine-d5	3.93	84	467095	31.84	ng/ul	0.00
7) Phenol-d5	7.42	99	574985	34.01	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.59	67	364228	34.67	ng/ul	0.00
11) 2-Chlorophenol-d4	7.80	132	485121	34.45	ng/ul	0.00
15) 4-Methylphenol-d8	8.99	113	456032	34.20	ng/ul	0.00
21) Nitrobenzene-d5	9.46	128	225927	34.51	ng/ul	0.00
24) 2-Nitrophenol-d4	10.19	143	241347	36.90	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.73	165	436718	32.92	ng/ul	0.00
31) 4-Chloroaniline-d4	11.25	131	640189	35.28	ng/ul	0.00
46) Dimethylphthalate-d6	14.30	166	1356076	37.41	ng/ul	0.00
49) Acenaphthylene-d8	14.61	160	1547715	33.26	ng/ul	0.00
54) 4-Nitrophenol-d4	15.07	143	226404	32.39	ng/ul	0.00
60) Fluorene-d10	15.89	176	1112278	34.24	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.99	200	169276	29.17	ng/ul	0.00
73) Anthracene-d10	17.73	188	1693362	34.55	ng/ul	0.00
81) Pyrene-d10	19.98	212	1774409	37.63	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.99	264	1638013	36.24	ng/ul	0.00

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

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