

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051021\
 Data File : BM030003.D
 Acq On : 13 May 2021 08:36
 Operator : CG/JU
 Sample : PB136264BL
 Misc :
 ALS Vial : 103 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK264

Quant Time: May 13 10:26:28 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050721.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 11 15:21:38 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	170228	20.00	ng/ul	-0.01
20) Naphthalene-d8	10.31	136	743346	20.00	ng/ul	-0.02
38) Acenaphthene-d10	14.19	164	483904	20.00	ng/ul	-0.02
64) Phenanthrene-d10	16.93	188	917929	20.00	ng/ul	-0.01
79) Chrysene-d12	21.13	240	679307	20.00	ng/ul	0.00
88) Perylene-d12	23.28	264	665575	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	29059	6.82	ng/uL	0.00
4) Pyridine-d5	3.47	84	333334	31.23	ng/ul	-0.01
7) Phenol-d5	6.73	99	484552	30.61	ng/ul	-0.01
9) Bis-(2-Chloroethyl)ether-d	6.89	67	320191	32.62	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.07	132	414272	33.91	ng/ul	-0.01
15) 4-Methylphenol-d8	8.26	113	401836	30.69	ng/ul	-0.01
21) Nitrobenzene-d5	8.69	128	195577	38.18	ng/ul	-0.02
24) 2-Nitrophenol-d4	9.41	143	220023	38.05	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	9.95	165	379678	32.75	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.46	131	543195	30.79	ng/ul	-0.01
46) Dimethylphthalate-d6	13.61	166	1368356	36.48	ng/ul	-0.01
49) Acenaphthylene-d8	13.88	160	1637540	34.45	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.43	143	134421	22.00	ng/ul	0.00
60) Fluorene-d10	15.19	176	1122287	35.25	ng/ul	-0.01
65) 4,6-Dinitro-2-methylphenol	15.33	200	168033	28.18	ng/ul	0.00
73) Anthracene-d10	17.03	188	1694623	36.66	ng/ul	-0.01
81) Pyrene-d10	19.33	212	1671736	47.97	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.14	264	1348278	35.74	ng/ul	0.00

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

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