

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111721\
 Data File : BM033143.D
 Acq On : 18 Nov 2021 05:58
 Operator : CG/JU
 Sample : PB140810BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK810

Quant Time: Nov 18 06:46:51 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM111721.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 17 14:14:11 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.972	152	116448	20.000	ng/u1	0.00
20) Naphthalene-d8	10.766	136	478797	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.589	164	319237	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.325	188	636501	20.000	ng/u1	0.00
79) Chrysene-d12	21.477	240	561369	20.000	ng/u1	0.00
88) Perylene-d12	23.824	264	535711	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.420	96	17672	5.851	ng/uL	0.00
4) Pyridine-d5	3.837	84	249585	29.997	ng/u1	-0.01
7) Phenol-d5	7.119	99	303756	30.838	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.296	67	199245	31.891	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.502	132	236111	31.648	ng/u1	0.00
15) 4-Methylphenol-d8	8.660	113	242030	31.661	ng/u1	-0.01
21) Nitrobenzene-d5	9.131	128	116076	33.767	ng/u1	0.00
24) 2-Nitrophenol-d4	9.849	143	122383	35.548	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.378	165	237184	30.084	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.901	131	336134	32.059	ng/u1	0.00
46) Dimethylphthalate-d6	13.995	166	756306	32.299	ng/u1	-0.01
49) Acenaphthylene-d8	14.284	160	933839	30.957	ng/u1	0.00
54) 4-Nitrophenol-d4	14.766	143	103439	27.120	ng/u1	0.00
60) Fluorene-d10	15.578	176	656967	31.276	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.683	200	83569	27.937	ng/u1	-0.01
73) Anthracene-d10	17.419	188	983108	32.042	ng/u1	-0.01
81) Pyrene-d10	19.701	212	1081241	32.593	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.671	264	915108	31.829	ng/u1	-0.01

Target Compounds Qvalue

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

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