

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110222\
 Data File : BM037321.D
 Acq On : 02 Nov 2022 18:05
 Operator : CG/JU
 Sample : PB148670BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK670

Quant Time: Nov 02 22:54:58 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 02 03:33:37 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.857	152	316324	20.000 ng/u1	0.00
20) Naphthalene-d8	10.657	136	1392973	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.492	164	900414	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.233	188	1851860	20.000 ng/u1	0.00
79) Chrysene-d12	21.403	240	1742498	20.000 ng/u1	0.00
88) Perylene-d12	23.756	264	1684209	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.275	96	49872	6.950 ng/uL	0.00
4) Pyridine-d5	3.704	84	171441	8.003 ng/u1	0.00
7) Phenol-d5	7.028	99	831938	30.220 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.192	67	546481	32.350 ng/u1	0.00
11) 2-Chlorophenol-d4	7.387	132	688519	32.735 ng/u1	0.00
15) 4-Methylphenol-d8	8.575	113	664566	29.563 ng/u1	0.00
21) Nitrobenzene-d5	9.028	128	338772	32.558 ng/u1	0.00
24) 2-Nitrophenol-d4	9.745	143	363418	32.063 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	628358	30.070 ng/u1	0.00
31) 4-Chloroaniline-d4	10.810	131	102156	3.199 ng/u1	0.00
46) Dimethylphthalate-d6	13.910	166	2065513	33.177 ng/u1	0.00
49) Acenaphthylene-d8	14.186	160	2429967	32.466 ng/u1	0.00
54) 4-Nitrophenol-d4	14.704	143	342912	27.556 ng/u1	0.00
60) Fluorene-d10	15.480	176	1919421	34.573 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.610	200	306141	26.101 ng/u1	0.00
73) Anthracene-d10	17.327	188	2998615	35.159 ng/u1	0.00
81) Pyrene-d10	19.621	212	3286900	36.098 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.603	264	2931905	33.818 ng/u1	0.00

Target Compounds Qvalue

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

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