

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
 Data File : BM036557.D
 Acq On : 16 Sep 2022 10:46
 Operator : CG/JU
 Sample : PB147637BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK637

Quant Time: Sep 16 22:58:06 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 16 02:39:54 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	8.022	152	303810	20.000 ng/u1	0.00
20) Naphthalene-d8	10.839	136	1336399	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.651	164	904349	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.386	188	1925165	20.000 ng/u1	0.00
79) Chrysene-d12	21.550	240	1781332	20.000 ng/u1	0.00
88) Perylene-d12	23.997	264	1734693	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.393	96	49013	6.048 ng/uL	0.00
4) Pyridine-d5	3.810	84	611441	25.870 ng/u1	0.00
7) Phenol-d5	7.169	99	770483	28.379 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	514528	31.206 ng/u1	0.00
11) 2-Chlorophenol-d4	7.551	132	624037	31.926 ng/u1	0.00
15) 4-Methylphenol-d8	8.722	113	606317	30.224 ng/u1	-0.01
21) Nitrobenzene-d5	9.187	128	291552	32.232 ng/u1	0.00
24) 2-Nitrophenol-d4	9.916	143	267738	33.045 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	549055	29.241 ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	943718	30.229 ng/u1	0.00
46) Dimethylphthalate-d6	14.057	166	2110502	34.602 ng/u1	0.00
49) Acenaphthylene-d8	14.345	160	2330231	32.208 ng/u1	0.00
54) 4-Nitrophenol-d4	14.821	143	308488	26.076 ng/u1	0.00
60) Fluorene-d10	15.639	176	1840066	33.664 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	184831	24.266 ng/u1	0.00
73) Anthracene-d10	17.480	188	2926343	33.245 ng/u1	0.00
81) Pyrene-d10	19.762	212	3265664	37.265 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.839	264	2997100	35.364 ng/u1	0.00

Target Compounds Qvalue

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

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