

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
 Data File : BM036635.D
 Acq On : 21 Sep 2022 21:48
 Operator : CG/JU
 Sample : PB147764BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK764

Quant Time: Sep 21 22:54:12 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 21 22:42:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.022	152	326506	20.000	ng/u1	0.00
20) Naphthalene-d8	10.833	136	1488832	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.645	164	996586	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.380	188	2063010	20.000	ng/u1	0.00
79) Chrysene-d12	21.544	240	1720996	20.000	ng/u1	0.00
88) Perylene-d12	23.985	264	1541800	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	44876	5.152	ng/uL	0.00
4) Pyridine-d5	3.822	84	495282	19.498	ng/u1	0.00
7) Phenol-d5	7.175	99	800780	27.445	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.345	67	500970	28.272	ng/u1	0.00
11) 2-Chlorophenol-d4	7.551	132	651290	31.004	ng/u1	0.00
15) 4-Methylphenol-d8	8.728	113	651973	30.240	ng/u1	0.00
21) Nitrobenzene-d5	9.186	128	325386	32.289	ng/u1	0.00
24) 2-Nitrophenol-d4	9.910	143	330424	36.607	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	606518	28.994	ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	593755	17.072	ng/u1	0.00
46) Dimethylphthalate-d6	14.057	166	2041461	30.372	ng/u1	0.00
49) Acenaphthylene-d8	14.339	160	2365814	29.674	ng/u1	0.00
54) 4-Nitrophenol-d4	14.833	143	339348	26.030	ng/u1	0.00
60) Fluorene-d10	15.633	176	1798129	29.852	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	209122	25.621	ng/u1	0.00
73) Anthracene-d10	17.480	188	2792395	29.604	ng/u1	0.00
81) Pyrene-d10	19.762	212	3031209	35.802	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.833	264	2358200	31.307	ng/u1	0.00

Target Compounds Qvalue

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

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