

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
 Data File : BM036651.D
 Acq On : 22 Sep 2022 07:57
 Operator : CG/JU
 Sample : PB147686BL
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK686

Quant Time: Sep 22 08:40:31 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	8.016	152	425344	20.000	ng/u1	0.00
20) Naphthalene-d8	10.834	136	1927628	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.645	164	1229950	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.380	188	2438798	20.000	ng/u1	0.00
79) Chrysene-d12	21.545	240	1745445	20.000	ng/u1	0.00
88) Perylene-d12	23.986	264	1527486	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	55359	4.879	ng/uL	0.00
4) Pyridine-d5	3.816	84	624264	18.865	ng/u1	0.00
7) Phenol-d5	7.175	99	1015905	26.727	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.346	67	635293	27.521	ng/u1	0.00
11) 2-Chlorophenol-d4	7.545	132	829762	30.321	ng/u1	0.00
15) 4-Methylphenol-d8	8.728	113	844044	30.052	ng/u1	0.00
21) Nitrobenzene-d5	9.187	128	419831	32.178	ng/u1	0.00
24) 2-Nitrophenol-d4	9.910	143	419753	35.917	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.445	165	786435	29.037	ng/u1	0.00
31) 4-Chloroaniline-d4	10.963	131	754984	16.766	ng/u1	0.00
46) Dimethylphthalate-d6	14.057	166	2495670	30.085	ng/u1	0.00
49) Acenaphthylene-d8	14.339	160	2980122	30.287	ng/u1	0.00
54) 4-Nitrophenol-d4	14.827	143	414457	25.759	ng/u1	0.00
60) Fluorene-d10	15.627	176	2233611	30.046	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.739	200	227584	23.586	ng/u1	0.00
73) Anthracene-d10	17.480	188	3373857	30.257	ng/u1	0.00
81) Pyrene-d10	19.762	212	3425350	39.890	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.827	264	2362904	31.663	ng/u1	0.00

Target Compounds Qvalue

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

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