

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092322\
 Data File : BM036710.D
 Acq On : 24 Sep 2022 12:07
 Operator : CG/JU
 Sample : PB147712BL
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK712

Quant Time: Sep 26 01:10:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 24 00:41:16 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.004	152	382359	20.000	ng/u1	0.00
20) Naphthalene-d8	10.816	136	1805196	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.633	164	1150500	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.368	188	2236357	20.000	ng/u1	0.00
79) Chrysene-d12	21.533	240	1527990	20.000	ng/u1	-0.01
88) Perylene-d12	23.968	264	1341475	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	58936	5.778	ng/uL	0.00
4) Pyridine-d5	3.810	84	838666	28.194	ng/u1	0.00
7) Phenol-d5	7.163	99	1115886	32.657	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.334	67	695817	33.532	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	918426	37.334	ng/u1	0.00
15) 4-Methylphenol-d8	8.716	113	922746	36.548	ng/u1	-0.01
21) Nitrobenzene-d5	9.175	128	458915	37.559	ng/u1	0.00
24) 2-Nitrophenol-d4	9.898	143	459884	42.020	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.433	165	849382	33.488	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.951	131	1360412	32.260	ng/u1	-0.01
46) Dimethylphthalate-d6	14.045	166	2657438	34.248	ng/u1	0.00
49) Acenaphthylene-d8	14.327	160	3222775	35.015	ng/u1	0.00
54) 4-Nitrophenol-d4	14.821	143	416846	27.697	ng/u1	-0.01
60) Fluorene-d10	15.621	176	2357986	33.910	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.733	200	206703	23.362	ng/u1	0.00
73) Anthracene-d10	17.468	188	3519193	34.417	ng/u1	0.00
81) Pyrene-d10	19.750	212	3515641	46.769	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.815	264	2402633	36.660	ng/u1	0.00

Target Compounds Qvalue

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

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