

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091922\
 Data File : BM036600.D
 Acq On : 20 Sep 2022 03:48
 Operator : CG/JU
 Sample : PB147643BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK643

Quant Time: Sep 20 04:24:10 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	8.028	152	548283	20.000 ng/u1	0.00
20) Naphthalene-d8	10.839	136	2527634	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.651	164	1668710	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.386	188	3387176	20.000 ng/u1	0.00
79) Chrysene-d12	21.550	240	3188653	20.000 ng/u1	0.00
88) Perylene-d12	23.991	264	2749864	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.393	96	71543	4.892 ng/uL	0.00
4) Pyridine-d5	3.816	84	975871	22.878 ng/u1	0.00
7) Phenol-d5	7.175	99	1314476	26.828 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	812255	27.297 ng/u1	0.00
11) 2-Chlorophenol-d4	7.551	132	1056173	29.941 ng/u1	0.00
15) 4-Methylphenol-d8	8.728	113	1078676	29.794 ng/u1	0.00
21) Nitrobenzene-d5	9.192	128	540761	31.608 ng/u1	0.00
24) 2-Nitrophenol-d4	9.916	143	543691	35.479 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	996517	28.060 ng/u1	0.00
31) 4-Chloroaniline-d4	10.969	131	1594890	27.011 ng/u1	0.00
46) Dimethylphthalate-d6	14.063	166	3286897	29.205 ng/u1	0.00
49) Acenaphthylene-d8	14.345	160	3940168	29.515 ng/u1	0.00
54) 4-Nitrophenol-d4	14.827	143	605120	27.721 ng/u1	0.00
60) Fluorene-d10	15.639	176	2913935	28.891 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	385933	28.798 ng/u1	0.00
73) Anthracene-d10	17.486	188	4615792	29.804 ng/u1	0.00
81) Pyrene-d10	19.762	212	5160391	32.896 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.838	264	4267596	31.766 ng/u1	0.00

Target Compounds Qvalue

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

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