

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
 Data File : BM036741.D
 Acq On : 26 Sep 2022 12:27
 Operator : CG/JU
 Sample : PB147828BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK828

Quant Time: Sep 26 23:12:37 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 26 23:10:22 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.004	152	373837	20.000	ng/u1	0.00
20) Naphthalene-d8	10.816	136	1610511	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.627	164	1028488	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.368	188	2095076	20.000	ng/u1	0.00
79) Chrysene-d12	21.533	240	1610977	20.000	ng/u1	0.00
88) Perylene-d12	23.968	264	1423604	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	50660	5.080	ng/uL	0.00
4) Pyridine-d5	3.810	84	668934	23.001	ng/u1	0.00
7) Phenol-d5	7.163	99	924936	27.686	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	589194	29.041	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	734288	30.529	ng/u1	0.00
15) 4-Methylphenol-d8	8.710	113	741385	30.034	ng/u1	0.00
21) Nitrobenzene-d5	9.169	128	359673	32.995	ng/u1	0.00
24) 2-Nitrophenol-d4	9.892	143	347415	35.581	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.433	165	670358	29.625	ng/u1	0.00
31) 4-Chloroaniline-d4	10.951	131	1001697	26.625	ng/u1	0.00
46) Dimethylphthalate-d6	14.039	166	2135207	30.782	ng/u1	0.00
49) Acenaphthylene-d8	14.327	160	2514944	30.566	ng/u1	0.00
54) 4-Nitrophenol-d4	14.815	143	326695	24.282	ng/u1	0.00
60) Fluorene-d10	15.615	176	1924455	30.958	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.727	200	183406	22.126	ng/u1	0.00
73) Anthracene-d10	17.462	188	2946893	30.763	ng/u1	0.00
81) Pyrene-d10	19.745	212	3126920	39.455	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.809	264	2184720	31.412	ng/u1	0.00

Target Compounds Qvalue

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

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