

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092322\
 Data File : BM036689.D
 Acq On : 23 Sep 2022 23:00
 Operator : CG/JU
 Sample : PB147828BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK828

Quant Time: Sep 24 00:51:03 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	8.010	152	390764	20.000	ng/ul	0.00
20) Naphthalene-d8	10.822	136	1791104	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.633	164	1162461	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.374	188	2317897	20.000	ng/ul	0.00
79) Chrysene-d12	21.539	240	1758131	20.000	ng/ul	0.00
88) Perylene-d12	23.980	264	1531813	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	54852	5.262	ng/uL	0.00
4) Pyridine-d5	3.817	84	692029	22.764	ng/ul	0.00
7) Phenol-d5	7.169	99	969771	27.771	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	598798	28.236	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	778386	30.961	ng/ul	0.00
15) 4-Methylphenol-d8	8.722	113	781392	30.283	ng/ul	0.00
21) Nitrobenzene-d5	9.175	128	385499	31.799	ng/ul	0.00
24) 2-Nitrophenol-d4	9.904	143	390151	35.929	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	723947	28.767	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	1078989	25.788	ng/ul	0.00
46) Dimethylphthalate-d6	14.051	166	2339955	29.846	ng/ul	0.00
49) Acenaphthylene-d8	14.333	160	2732457	29.382	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	371259	24.414	ng/ul	0.00
60) Fluorene-d10	15.622	176	2063416	29.368	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.733	200	192092	20.946	ng/ul	0.00
73) Anthracene-d10	17.474	188	3082529	29.086	ng/ul	0.00
81) Pyrene-d10	19.757	212	3240557	37.466	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264	2315239	30.937	ng/ul	0.00

Target Compounds Qvalue

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

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