

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
 Data File : BM036752.D
 Acq On : 26 Sep 2022 19:21
 Operator : CG/JU
 Sample : PB147885BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK885

Quant Time: Sep 26 23:15:36 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	338126	20.000	ng/u1	0.00
20) Naphthalene-d8	10.816	136	1464948	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.627	164	911671	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.362	188	1775845	20.000	ng/u1	0.00
79) Chrysene-d12	21.533	240	1260077	20.000	ng/u1	0.00
88) Perylene-d12	23.968	264	1070109	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	60694	6.729	ng/uL	0.00
4) Pyridine-d5	3.810	84	774561	29.445	ng/u1	0.00
7) Phenol-d5	7.163	99	1070829	35.438	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	674382	36.750	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	844445	38.817	ng/u1	0.00
15) 4-Methylphenol-d8	8.710	113	860368	38.535	ng/u1	0.00
21) Nitrobenzene-d5	9.169	128	424933	42.855	ng/u1	0.00
24) 2-Nitrophenol-d4	9.892	143	410563	46.226	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.433	165	772258	37.519	ng/u1	0.00
31) 4-Chloroaniline-d4	10.951	131	1156775	33.803	ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	2466319	40.111	ng/u1	0.00
49) Acenaphthylene-d8	14.327	160	2906569	39.852	ng/u1	0.00
54) 4-Nitrophenol-d4	14.815	143	354407	29.717	ng/u1	0.00
60) Fluorene-d10	15.615	176	2176314	39.496	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.727	200	206340	29.368	ng/u1	0.00
73) Anthracene-d10	17.462	188	3145980	38.745	ng/u1	0.00
81) Pyrene-d10	19.750	212	3225145	52.026	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.809	264	2140997	40.952	ng/u1	0.00

Target Compounds Qvalue

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

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