

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
 Data File : BM036773.D
 Acq On : 27 Sep 2022 09:29
 Operator : CG/JU
 Sample : PB147764BL
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK764

Quant Time: Sep 27 11:51:51 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	7.993	152	332811	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.804	136	1433010	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.622	164	862118	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.357	188	1579901	20.000	ng/u1	-0.01
79) Chrysene-d12	21.527	240	1075550	20.000	ng/u1	0.00
88) Perylene-d12	23.956	264	942169	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	55919	6.299	ng/uL	0.00
4) Pyridine-d5	3.811	84	621402	24.000	ng/u1	0.00
7) Phenol-d5	7.151	99	969036	32.582	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.322	67	611278	33.844	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.522	132	770511	35.984	ng/u1	-0.01
15) 4-Methylphenol-d8	8.704	113	772366	35.146	ng/u1	-0.01
21) Nitrobenzene-d5	9.163	128	378427	39.016	ng/u1	0.00
24) 2-Nitrophenol-d4	9.887	143	365007	42.013	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.422	165	690376	34.289	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.939	131	391207	11.686	ng/u1	-0.01
46) Dimethylphthalate-d6	14.033	166	2107261	36.241	ng/u1	0.00
49) Acenaphthylene-d8	14.316	160	2480229	35.961	ng/u1	0.00
54) 4-Nitrophenol-d4	14.810	143	281699	24.978	ng/u1	-0.01
60) Fluorene-d10	15.610	176	1828674	35.095	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	165287	26.443	ng/u1	-0.01
73) Anthracene-d10	17.457	188	2604306	36.052	ng/u1	0.00
81) Pyrene-d10	19.739	212	2553135	48.252	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.797	264	1739131	37.782	ng/u1	0.00

Target Compounds Qvalue

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

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