

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092622\
 Data File : BM036770.D
 Acq On : 27 Sep 2022 06:52
 Operator : CG/JU
 Sample : PB147886TB 10X
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLEB886

Quant Time: Sep 27 07:37:21 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	7.998	152	422790	20.000	ng/u1	0.00
20) Naphthalene-d8	10.804	136	1833539	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.621	164	1115756	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.362	188	2160053	20.000	ng/u1	0.00
79) Chrysene-d12	21.527	240	1554903	20.000	ng/u1	0.00
88) Perylene-d12	23.956	264	1325729	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	67549	5.989	ng/uL	0.00
4) Pyridine-d5	3.810	84	901301	27.402	ng/u1	0.00
7) Phenol-d5	7.157	99	1251093	33.113	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.328	67	782024	34.082	ng/u1	0.00
11) 2-Chlorophenol-d4	7.528	132	996706	36.641	ng/u1	0.00
15) 4-Methylphenol-d8	8.704	113	1000002	35.820	ng/u1	-0.01
21) Nitrobenzene-d5	9.163	128	499234	40.227	ng/u1	0.00
24) 2-Nitrophenol-d4	9.886	143	487227	43.830	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.422	165	899240	34.906	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.945	131	1316880	30.745	ng/u1	0.00
46) Dimethylphthalate-d6	14.039	166	2718436	36.125	ng/u1	0.00
49) Acenaphthylene-d8	14.316	160	3277745	36.721	ng/u1	0.00
54) 4-Nitrophenol-d4	14.810	143	412640	28.271	ng/u1	-0.01
60) Fluorene-d10	15.610	176	2428298	36.008	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.721	200	245290	28.702	ng/u1	-0.01
73) Anthracene-d10	17.457	188	3575487	36.203	ng/u1	0.00
81) Pyrene-d10	19.745	212	3663843	47.896	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.803	264	2413979	37.270	ng/u1	0.00

Target Compounds Qvalue

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

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