

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
 Data File : BM047750.D
 Acq On : 01 Oct 2024 20:19
 Operator : RC/JU
 Sample : PB163513BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK513

Quant Time: Oct 01 23:18:10 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 28 02:30:36 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	263427	20.000	ng/u1	0.00
20) Naphthalene-d8	10.604	136	1066210	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.445	164	659044	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.180	188	1338772	20.000	ng/u1	0.00
79) Chrysene-d12	21.403	240	1241551	20.000	ng/u1	0.00
88) Perylene-d12	24.409	264	1467098	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	31053	3.809	ng/uL	0.00
4) Pyridine-d5	3.669	84	453037	18.981	ng/u1	0.00
7) Phenol-d5	6.975	99	473498	19.009	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	305488	17.283	ng/u1	0.00
11) 2-Chlorophenol-d4	7.340	132	402286	22.197	ng/u1	0.00
15) 4-Methylphenol-d8	8.510	113	369226	19.936	ng/u1	0.00
21) Nitrobenzene-d5	8.963	128	187933	22.138	ng/u1	0.00
24) 2-Nitrophenol-d4	9.687	143	188999	23.232	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	353803	21.771	ng/u1	0.00
31) 4-Chloroaniline-d4	10.734	131	533123	21.219	ng/u1	0.00
46) Dimethylphthalate-d6	13.851	166	1014776	21.238	ng/u1	0.00
49) Acenaphthylene-d8	14.133	160	1210455	21.366	ng/u1	0.00
54) 4-Nitrophenol-d4	14.633	143	158974	16.711	ng/u1	0.00
60) Fluorene-d10	15.433	176	892289	21.560	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.545	200	103880	13.260	ng/u1	0.00
73) Anthracene-d10	17.274	188	1356817	20.611	ng/u1	-0.01
81) Pyrene-d10	19.562	212	1542797	21.895	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.203	264	1576401	20.710	ng/u1	0.00

Target Compounds Qvalue

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

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