

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122123\
 Data File : BM043532.D
 Acq On : 22 Dec 2023 10:17
 Operator : MA/JU
 Sample : PB157967BL
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK967

Quant Time: Dec 23 01:27:01 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 21 21:59:32 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	234343	20.000	ng/ul	0.00
20) Naphthalene-d8	10.775	136	985794	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.604	164	536931	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.351	188	1072253	20.000	ng/ul	0.00
79) Chrysene-d12	21.533	240	879565	20.000	ng/ul	0.00
88) Perylene-d12	23.950	264	1034177	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	35366	4.905	ng/uL	0.00
4) Pyridine-d5	3.787	84	499484	24.069	ng/ul	0.00
7) Phenol-d5	7.122	99	574962	21.568	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.304	67	382265	22.642	ng/ul	0.00
11) 2-Chlorophenol-d4	7.492	132	431986	21.223	ng/ul	0.00
15) 4-Methylphenol-d8	8.675	113	420477	20.326	ng/ul	0.00
21) Nitrobenzene-d5	9.139	128	205517	21.548	ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	198686	20.161	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.392	165	349117	20.068	ng/ul	0.00
31) 4-Chloroaniline-d4	10.922	131	592654	23.157	ng/ul	0.00
46) Dimethylphthalate-d6	14.021	166	1050306	21.730	ng/ul	0.00
49) Acenaphthylene-d8	14.298	160	1259096	21.959	ng/ul	0.00
54) 4-Nitrophenol-d4	14.804	143	186411	20.377	ng/ul	0.00
60) Fluorene-d10	15.598	176	926812	23.188	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.715	200	101942	14.375	ng/ul	0.00
73) Anthracene-d10	17.451	188	1406837	23.395	ng/ul	0.00
81) Pyrene-d10	19.739	212	1473435	21.218	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.791	264	1500197	23.507	ng/ul	0.00

Target Compounds Qvalue

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

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