

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081222\
 Data File : BM036232.D
 Acq On : 12 Aug 2022 16:17
 Operator : CG/JU
 Sample : PB146934BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK934

Quant Time: Aug 12 17:22:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM072522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 12 11:23:11 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	190484	20.000	ng/u1	0.00
20) Naphthalene-d8	10.622	136	783791	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.457	164	445933	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.203	188	878599	20.000	ng/u1	0.00
79) Chrysene-d12	21.386	240	846667	20.000	ng/u1	0.00
88) Perylene-d12	23.727	264	927451	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	30709	6.066	ng/uL	0.00
4) Pyridine-d5	3.646	84	310377	22.381	ng/u1	0.00
7) Phenol-d5	7.004	99	496058	31.054	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.151	67	306139	28.749	ng/u1	0.00
11) 2-Chlorophenol-d4	7.351	132	415172	31.312	ng/u1	0.00
15) 4-Methylphenol-d8	8.551	113	374137	28.912	ng/u1	0.00
21) Nitrobenzene-d5	8.986	128	201095	32.378	ng/u1	0.00
24) 2-Nitrophenol-d4	9.710	143	208840	34.513	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.257	165	342683	31.312	ng/u1	0.00
31) 4-Chloroaniline-d4	10.769	131	420648	22.705	ng/u1	0.00
46) Dimethylphthalate-d6	13.880	166	995961	32.788	ng/u1	0.00
49) Acenaphthylene-d8	14.157	160	1208805	31.190	ng/u1	0.00
54) 4-Nitrophenol-d4	14.698	143	158036	25.336	ng/u1	0.00
60) Fluorene-d10	15.451	176	899206	32.759	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.580	200	126590	26.804	ng/u1	0.00
73) Anthracene-d10	17.303	188	1390749	35.012	ng/u1	0.00
81) Pyrene-d10	19.597	212	1519345	32.379	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.574	264	1635850	34.742	ng/u1	0.00

Target Compounds Qvalue

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

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