

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081222\
 Data File : BM036259.D
 Acq On : 13 Aug 2022 09:45
 Operator : CG/JU
 Sample : PB146949BL
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK949

Quant Time: Aug 14 23:00:24 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	7.810	152	220781	20.000	ng/u1	0.00
20) Naphthalene-d8	10.610	136	950774	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.451	164	564472	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.198	188	1171786	20.000	ng/u1	0.00
79) Chrysene-d12	21.380	240	1187132	20.000	ng/u1	0.00
88) Perylene-d12	23.715	264	1278381	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.216	96	31361	5.345	ng/uL	0.00
4) Pyridine-d5	3.640	84	325452	20.247	ng/u1	0.00
7) Phenol-d5	6.998	99	518165	27.987	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.145	67	329617	26.706	ng/u1	0.00
11) 2-Chlorophenol-d4	7.345	132	427392	27.810	ng/u1	0.00
15) 4-Methylphenol-d8	8.539	113	398270	26.553	ng/u1	-0.01
21) Nitrobenzene-d5	8.981	128	205920	27.332	ng/u1	0.00
24) 2-Nitrophenol-d4	9.698	143	214068	29.164	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.245	165	355552	26.782	ng/u1	0.00
31) 4-Chloroaniline-d4	10.763	131	446382	19.862	ng/u1	0.00
46) Dimethylphthalate-d6	13.874	166	1123871	29.229	ng/u1	0.00
49) Acenaphthylene-d8	14.145	160	1320522	26.917	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.686	143	193730	24.536	ng/u1	-0.01
60) Fluorene-d10	15.445	176	1015314	29.221	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.574	200	146905	23.323	ng/u1	0.00
73) Anthracene-d10	17.292	188	1620530	30.589	ng/u1	-0.01
81) Pyrene-d10	19.586	212	1857202	28.228	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.562	264	1971804	30.381	ng/u1	-0.01

Target Compounds Qvalue

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

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