

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112522\
 Data File : BM037742.D
 Acq On : 26 Nov 2022 20:33
 Operator : CG/JU
 Sample : PB149109BL
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK109

Quant Time: Nov 27 21:30:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 24 01:27:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.110	152	282658	20.000	ng/ul	0.00
20) Naphthalene-d8	10.933	136	1273299	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.733	164	854932	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.468	188	1779617	20.000	ng/ul	0.00
79) Chrysene-d12	21.633	240	1636232	20.000	ng/ul	0.00
88) Perylene-d12	24.127	264	1405364	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	53022	7.779	ng/uL	0.00
4) Pyridine-d5	3.875	84	566629	25.541	ng/ul	0.00
7) Phenol-d5	7.257	99	749174	26.567	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.428	67	492151	26.163	ng/ul	0.00
11) 2-Chlorophenol-d4	7.634	132	569402	29.059	ng/ul	0.00
15) 4-Methylphenol-d8	8.816	113	597840	26.642	ng/ul	0.00
21) Nitrobenzene-d5	9.280	128	310562	31.112	ng/ul	0.00
24) 2-Nitrophenol-d4	10.004	143	332133	33.768	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.545	165	560546	31.142	ng/ul	0.00
31) 4-Chloroaniline-d4	11.069	131	861572	28.846	ng/ul	0.00
46) Dimethylphthalate-d6	14.139	166	1900513	31.384	ng/ul	0.00
49) Acenaphthylene-d8	14.433	160	2186715	30.593	ng/ul	0.00
54) 4-Nitrophenol-d4	14.915	143	364496	27.326	ng/ul	0.00
60) Fluorene-d10	15.721	176	1661470	31.327	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.827	200	266277	27.815	ng/ul	-0.01
73) Anthracene-d10	17.568	188	2685265	32.267	ng/ul	0.00
81) Pyrene-d10	19.850	212	3021649	36.340	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.968	264	2399832	33.175	ng/ul	0.00

Target Compounds Qvalue

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

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