

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112522\
 Data File : BM037732.D
 Acq On : 26 Nov 2022 14:17
 Operator : CG/JU
 Sample : PB149109BL
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK109

Quant Time: Nov 27 21:28:02 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	207964	20.000	ng/u1	0.00
20) Naphthalene-d8	10.933	136	1020430	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.739	164	641196	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.474	188	1287440	20.000	ng/u1	0.00
79) Chrysene-d12	21.633	240	1137013	20.000	ng/u1	0.00
88) Perylene-d12	24.127	264	1039791	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	42214	8.418	ng/uL	0.00
4) Pyridine-d5	3.869	84	482011	29.530	ng/u1	0.00
7) Phenol-d5	7.257	99	606463	29.231	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.428	67	428366	30.951	ng/u1	0.00
11) 2-Chlorophenol-d4	7.633	132	445322	30.889	ng/u1	0.00
15) 4-Methylphenol-d8	8.816	113	474156	28.719	ng/u1	0.00
21) Nitrobenzene-d5	9.280	128	245286	30.662	ng/u1	0.00
24) 2-Nitrophenol-d4	10.010	143	238797	30.295	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.545	165	407001	28.215	ng/u1	0.00
31) 4-Chloroaniline-d4	11.069	131	692499	28.930	ng/u1	0.00
46) Dimethylphthalate-d6	14.139	166	1429701	31.479	ng/u1	0.00
49) Acenaphthylene-d8	14.433	160	1645954	30.704	ng/u1	0.00
54) 4-Nitrophenol-d4	14.915	143	275436	27.532	ng/u1	0.00
60) Fluorene-d10	15.721	176	1219518	30.659	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.833	200	159580	23.042	ng/u1	0.00
73) Anthracene-d10	17.574	188	1918078	31.859	ng/u1	0.00
81) Pyrene-d10	19.850	212	2060746	35.665	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.968	264	1745886	32.621	ng/u1	0.00

Target Compounds Qvalue

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

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