

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110322\
 Data File : BM037363.D
 Acq On : 04 Nov 2022 01:34
 Operator : CG/JU
 Sample : PB148648BL
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK648

Quant Time: Nov 04 02:41:56 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 04 02:39:09 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	321487	20.000	ng/u1	0.00
20) Naphthalene-d8	10.645	136	1461233	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.480	164	961999	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.221	188	1963833	20.000	ng/u1	0.00
79) Chrysene-d12	21.392	240	1735555	20.000	ng/u1	0.00
88) Perylene-d12	23.739	264	1471220	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	47958	6.576	ng/uL	0.00
4) Pyridine-d5	3.693	84	679918	31.231	ng/u1	0.00
7) Phenol-d5	7.022	99	875648	31.297	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	564945	32.906	ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	721436	33.749	ng/u1	0.00
15) 4-Methylphenol-d8	8.563	113	714910	31.292	ng/u1	0.00
21) Nitrobenzene-d5	9.016	128	362166	33.180	ng/u1	0.00
24) 2-Nitrophenol-d4	9.734	143	389936	32.796	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	686096	31.299	ng/u1	0.00
31) 4-Chloroaniline-d4	10.798	131	1028269	30.696	ng/u1	0.00
46) Dimethylphthalate-d6	13.904	166	2168107	32.596	ng/u1	0.00
49) Acenaphthylene-d8	14.175	160	2657092	33.228	ng/u1	0.00
54) 4-Nitrophenol-d4	14.698	143	343857	25.863	ng/u1	0.00
60) Fluorene-d10	15.475	176	2051620	34.588	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	299808	24.103	ng/u1	0.00
73) Anthracene-d10	17.321	188	3212624	35.521	ng/u1	0.00
81) Pyrene-d10	19.610	212	3511108	38.714	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.586	264	2742704	36.216	ng/u1	0.00

Target Compounds Qvalue

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

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