

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110222\
 Data File : BM037311.D
 Acq On : 02 Nov 2022 10:47
 Operator : CG/JU
 Sample : PB148668BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK668

Quant Time: Nov 02 22:52:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 02 03:33:37 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.851	152	304245	20.000 ng/u1	0.00
20) Naphthalene-d8	10.657	136	1328538	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.486	164	856837	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.233	188	1783547	20.000 ng/u1	0.00
79) Chrysene-d12	21.403	240	1633747	20.000 ng/u1	0.00
88) Perylene-d12	23.750	264	1463144	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.275	96	40993	5.939 ng/uL	0.00
4) Pyridine-d5	3.693	84	561111	27.235 ng/u1	0.00
7) Phenol-d5	7.028	99	695063	26.251 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.192	67	462866	28.488 ng/u1	0.00
11) 2-Chlorophenol-d4	7.387	132	575319	28.439 ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	561258	25.958 ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	278972	28.111 ng/u1	0.00
24) 2-Nitrophenol-d4	9.745	143	301869	27.925 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	529223	26.554 ng/u1	0.00
31) 4-Chloroaniline-d4	10.804	131	813779	26.720 ng/u1	0.00
46) Dimethylphthalate-d6	13.910	166	1793707	30.277 ng/u1	0.00
49) Acenaphthylene-d8	14.186	160	2083302	29.250 ng/u1	0.00
54) 4-Nitrophenol-d4	14.704	143	272090	22.977 ng/u1	0.00
60) Fluorene-d10	15.480	176	1644081	31.120 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.610	200	245568	21.738 ng/u1	0.00
73) Anthracene-d10	17.327	188	2596866	31.615 ng/u1	0.00
81) Pyrene-d10	19.615	212	2845912	33.335 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.597	264	2410254	32.002 ng/u1	0.00

Target Compounds Qvalue

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

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