

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110322\
 Data File : BM037347.D
 Acq On : 03 Nov 2022 15:25
 Operator : CG/JU
 Sample : PB148651BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK651

Quant Time: Nov 04 01:03:29 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	299714	20.000	ng/u1	0.00
20) Naphthalene-d8	10.651	136	1334550	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.486	164	872149	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.227	188	1841577	20.000	ng/u1	0.00
79) Chrysene-d12	21.398	240	1807661	20.000	ng/u1	0.00
88) Perylene-d12	23.744	264	1662617	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	44119	6.489	ng/uL	0.00
4) Pyridine-d5	3.693	84	615978	30.350	ng/u1	0.00
7) Phenol-d5	7.022	99	796076	30.520	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	515360	32.198	ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	659992	33.117	ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	645238	30.294	ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	328516	32.954	ng/u1	0.00
24) 2-Nitrophenol-d4	9.739	143	355271	32.717	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	623893	31.163	ng/u1	0.00
31) 4-Chloroaniline-d4	10.798	131	940003	30.725	ng/u1	0.00
46) Dimethylphthalate-d6	13.904	166	1961353	32.525	ng/u1	0.00
49) Acenaphthylene-d8	14.180	160	2358154	32.528	ng/u1	0.00
54) 4-Nitrophenol-d4	14.704	143	332956	27.623	ng/u1	0.00
60) Fluorene-d10	15.480	176	1849475	34.393	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	283471	24.303	ng/u1	0.00
73) Anthracene-d10	17.327	188	2955734	34.850	ng/u1	0.00
81) Pyrene-d10	19.615	212	3360922	35.580	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.597	264	2999994	35.053	ng/u1	0.00

Target Compounds	Qvalue

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

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