

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110322\
 Data File : BM037339.D
 Acq On : 03 Nov 2022 09:52
 Operator : CG/JU
 Sample : PB148708BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK708

Quant Time: Nov 04 00:59:44 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	288365	20.000 ng/u1	0.00
20) Naphthalene-d8	10.651	136	1311145	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.486	164	878099	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.227	188	1920842	20.000 ng/u1	0.00
79) Chrysene-d12	21.398	240	1944091	20.000 ng/u1	0.00
88) Perylene-d12	23.739	264	1708299	20.000 ng/u1	-0.02
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.269	96	35987	5.501 ng/uL	0.00
4) Pyridine-d5	3.693	84	275425	14.104 ng/u1	0.00
7) Phenol-d5	7.022	99	627199	24.992 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	405891	26.357 ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	516906	26.958 ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	515377	25.149 ng/u1	0.00
21) Nitrobenzene-d5	9.016	128	259468	26.492 ng/u1	-0.01
24) 2-Nitrophenol-d4	9.740	143	280500	26.292 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	507133	25.783 ng/u1	0.00
31) 4-Chloroaniline-d4	10.798	131	406120	13.511 ng/u1	0.00
46) Dimethylphthalate-d6	13.904	166	1775659	29.247 ng/u1	0.00
49) Acenaphthylene-d8	14.181	160	2042711	27.986 ng/u1	0.00
54) 4-Nitrophenol-d4	14.698	143	279382	23.021 ng/u1	-0.01
60) Fluorene-d10	15.475	176	1653912	30.548 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	248116	20.394 ng/u1	0.00
73) Anthracene-d10	17.321	188	2666573	30.143 ng/u1	-0.01
81) Pyrene-d10	19.610	212	2959049	29.127 ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.592	264	2609159	29.671 ng/u1	-0.01

Target Compounds Qvalue

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

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