

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082222\
 Data File : BM036362.D
 Acq On : 22 Aug 2022 12:36
 Operator : CG/JU
 Sample : PB147153BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK153

Quant Time: Aug 23 01:00:03 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 23 00:55:01 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	236968	20.000	ng/u1	0.00
20) Naphthalene-d8	10.593	136	1028395	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.434	164	542345	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.180	188	919815	20.000	ng/u1	0.00
79) Chrysene-d12	21.362	240	765222	20.000	ng/u1	0.00
88) Perylene-d12	23.692	264	801303	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	34174	5.496	ng/uL	0.00
4) Pyridine-d5	3.628	84	511228	29.815	ng/u1	0.00
7) Phenol-d5	6.969	99	611674	30.240	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	421371	33.095	ng/u1	0.00
11) 2-Chlorophenol-d4	7.322	132	485807	30.568	ng/u1	0.00
15) 4-Methylphenol-d8	8.516	113	463504	30.690	ng/u1	0.00
21) Nitrobenzene-d5	8.963	128	244409	31.346	ng/u1	0.00
24) 2-Nitrophenol-d4	9.681	143	223464	28.255	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	391586	27.710	ng/u1	0.00
31) 4-Chloroaniline-d4	10.746	131	670391	30.600	ng/u1	0.00
46) Dimethylphthalate-d6	13.857	166	1139783	33.219	ng/u1	0.00
49) Acenaphthylene-d8	14.128	160	1400130	32.718	ng/u1	0.00
54) 4-Nitrophenol-d4	14.669	143	100992	15.534	ng/u1	0.00
60) Fluorene-d10	15.428	176	987036	31.835	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	71598	15.141	ng/u1	0.00
73) Anthracene-d10	17.275	188	1369375	33.481	ng/u1	0.00
81) Pyrene-d10	19.574	212	1401150	35.298	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.539	264	1361879	34.664	ng/u1	0.00

Target Compounds Qvalue

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

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