

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081922\
 Data File : BM036357.D
 Acq On : 19 Aug 2022 15:18
 Operator : CG/JU
 Sample : PB146669BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK669

Quant Time: Aug 20 02:19:58 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 16 15:00:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	291134	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.592	136	1261875	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.439	164	681436	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.180	188	1223001	20.000	ng/u1	-0.01
79) Chrysene-d12	21.368	240	1061808	20.000	ng/u1	0.00
88) Perylene-d12	23.697	264	1076940	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	38823	5.082	ng/uL	0.00
4) Pyridine-d5	3.628	84	552409	26.222	ng/u1	-0.01
7) Phenol-d5	6.969	99	706476	28.429	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.128	67	464100	29.669	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.322	132	580362	29.724	ng/u1	-0.01
15) 4-Methylphenol-d8	8.516	113	548528	29.562	ng/u1	-0.01
21) Nitrobenzene-d5	8.963	128	288921	30.199	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.681	143	299804	30.893	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.222	165	494218	28.501	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.739	131	763844	28.415	ng/u1	-0.02
46) Dimethylphthalate-d6	13.857	166	1374072	31.873	ng/u1	-0.01
49) Acenaphthylene-d8	14.133	160	1647405	30.638	ng/u1	0.00
54) 4-Nitrophenol-d4	14.669	143	189115	23.151	ng/u1	-0.01
60) Fluorene-d10	15.433	176	1186692	30.462	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	135882	21.611	ng/u1	0.00
73) Anthracene-d10	17.280	188	1688716	31.054	ng/u1	0.00
81) Pyrene-d10	19.574	212	1794481	32.579	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.544	264	1707202	32.332	ng/u1	0.00

Target Compounds Qvalue

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

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