

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081822\
 Data File : BM036344.D
 Acq On : 18 Aug 2022 12:16
 Operator : CG/JU
 Sample : PB146656BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK656

Quant Time: Aug 18 15:46:42 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	7.792	152	221481	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.592	136	961846	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.439	164	544792	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.180	188	1057346	20.000	ng/u1	-0.01
79) Chrysene-d12	21.368	240	977868	20.000	ng/u1	0.00
88) Perylene-d12	23.697	264	1019756	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.204	96	30201	5.197	ng/uL	0.00
4) Pyridine-d5	3.634	84	346592	21.627	ng/u1	0.00
7) Phenol-d5	6.975	99	511957	27.080	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.134	67	332134	27.910	ng/u1	0.00
11) 2-Chlorophenol-d4	7.328	132	430384	28.974	ng/u1	0.00
15) 4-Methylphenol-d8	8.516	113	398843	28.255	ng/u1	-0.01
21) Nitrobenzene-d5	8.963	128	209820	28.772	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.686	143	216674	29.292	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	361695	27.365	ng/u1	0.00
31) 4-Chloroaniline-d4	10.745	131	543446	26.522	ng/u1	-0.01
46) Dimethylphthalate-d6	13.857	166	1077406	31.260	ng/u1	-0.01
49) Acenaphthylene-d8	14.133	160	1353192	31.479	ng/u1	0.00
54) 4-Nitrophenol-d4	14.668	143	164274	25.154	ng/u1	-0.01
60) Fluorene-d10	15.433	176	948168	30.444	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	125938	23.168	ng/u1	0.00
73) Anthracene-d10	17.280	188	1465900	31.180	ng/u1	0.00
81) Pyrene-d10	19.574	212	1633503	32.203	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.550	264	1591307	31.827	ng/u1	0.00

Target Compounds Qvalue

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

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