

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060721\
 Data File : BM030306.D
 Acq On : 07 Jun 2021 10:34
 Operator : CG/JU
 Sample : PB136695BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK695

Quant Time: Jun 07 11:22:44 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM060621.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 07 09:55:51 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.851	152	191314	20.000	ng/ul	0.00
20) Naphthalene-d8	10.639	136	802070	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.474	164	539232	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.210	188	1147985	20.000	ng/ul	0.00
79) Chrysene-d12	21.368	240	1194669	20.000	ng/ul	0.00
88) Perylene-d12	23.656	264	1308604	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.316	96	30080	6.004	ng/uL	0.00
4) Pyridine-d5	3.728	84	353811	26.875	ng/ul	0.00
7) Phenol-d5	7.004	99	467290	28.103	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	311164	28.995	ng/ul	0.00
11) 2-Chlorophenol-d4	7.381	132	386299	32.124	ng/ul	0.00
15) 4-Methylphenol-d8	8.545	113	376566	28.357	ng/ul	0.00
21) Nitrobenzene-d5	9.004	128	179421	32.680	ng/ul	0.00
24) 2-Nitrophenol-d4	9.728	143	181832	33.346	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.257	165	371676	31.292	ng/ul	0.00
31) 4-Chloroaniline-d4	10.775	131	536657	29.555	ng/ul	0.00
46) Dimethylphthalate-d6	13.880	166	1281906	34.641	ng/ul	0.00
49) Acenaphthylene-d8	14.169	160	1514429	32.759	ng/ul	0.00
54) 4-Nitrophenol-d4	14.651	143	195426	28.214	ng/ul	0.00
60) Fluorene-d10	15.463	176	1122266	33.875	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.574	200	172064	25.966	ng/ul	0.00
73) Anthracene-d10	17.310	188	1792401	34.328	ng/ul	0.00
81) Pyrene-d10	19.592	212	2056875	32.894	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.509	264	2286059	35.034	ng/ul	0.00

Target Compounds Qvalue

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

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