

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060421\
 Data File : BM030285.D
 Acq On : 03 Jun 2021 16:28
 Operator : CG/JU
 Sample : PB136695BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK695

Quant Time: Jun 03 17:03:28 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM052621.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 03 16:38:49 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	201568	20.000	ng/ul	0.00
20) Naphthalene-d8	10.645	136	848945	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.480	164	521251	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.215	188	967803	20.000	ng/ul	0.00
79) Chrysene-d12	21.374	240	806382	20.000	ng/ul	0.00
88) Perylene-d12	23.656	264	829058	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.322	96	32413	6.581	ng/uL	0.00
4) Pyridine-d5	3.734	84	389656	28.220	ng/ul	0.00
7) Phenol-d5	7.010	99	517615	29.207	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	345389	30.875	ng/ul	0.00
11) 2-Chlorophenol-d4	7.387	132	413548	31.151	ng/ul	0.00
15) 4-Methylphenol-d8	8.551	113	404838	28.768	ng/ul	0.00
21) Nitrobenzene-d5	9.010	128	186280	33.617	ng/ul	0.00
24) 2-Nitrophenol-d4	9.733	143	182325	34.547	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.263	165	382536	28.092	ng/ul	0.00
31) 4-Chloroaniline-d4	10.780	131	556383	27.512	ng/ul	0.00
46) Dimethylphthalate-d6	13.886	166	1148541	28.821	ng/ul	0.00
49) Acenaphthylene-d8	14.174	160	1447039	28.170	ng/ul	0.00
54) 4-Nitrophenol-d4	14.657	143	159170	24.717	ng/ul	0.00
60) Fluorene-d10	15.468	176	1005156	28.896	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.580	200	122762	26.165	ng/ul	0.00
73) Anthracene-d10	17.309	188	1456062	30.398	ng/ul	0.00
81) Pyrene-d10	19.598	212	1483847	34.329	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.515	264	1407439	29.948	ng/ul	0.00

Target Compounds Qvalue

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

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