

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071621\
 Data File : BM030997.D
 Acq On : 16 Jul 2021 19:43
 Operator : CG/JU
 Sample : M2905-21DL 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBK32DL

Quant Time: Jul 16 23:19:13 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM071321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 16 16:49:56 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.649	152	1198	0.40	ng/ul	0.00
4) Naphthalene-d8	10.415	136	4585	0.40	ng/ul	# 0.00
9) Acenaphthene-d10	14.277	164	4803	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.027	188	7576	0.40	ng/ul	# 0.00
17) Chrysene-d12	21.212	240	6339	0.40	ng/ul	0.00
23) Perylene-d12	23.396	264	5922	0.40	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.224	96	671	0.54	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.007	152	770	0.10	ng/ul	0.00
18) Fluoranthene-d10	19.054	212	2029	0.10	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.258	88	270	0.20	ng/ul#	81
5) Naphthalene	10.464	128	33887	2.48	ng/ul	95
7) 2-Methylnaphthalene	12.083	142	15100	1.56	ng/ul	98
8) 1-Methylnaphthalene	12.304	142	10610	1.12	ng/ul	96
11) Acenaphthene	14.341	153	708	0.04	ng/ul#	84
12) Fluorene	15.331	166	1700	0.08	ng/ul#	94
14) Pentachlorophenol	16.676	266	65	0.02	ng/ul	93
15) Phenanthrene	17.065	178	1619	0.07	ng/ul#	30
19) Fluoranthene	19.058	202	652	0.02	ng/ul#	8

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

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