

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080221\
 Data File : BM031462.D
 Acq On : 05 Aug 2021 00:05
 Operator : CG/JU
 Sample : M3141-12
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGER0

Quant Time: Aug 05 01:02:33 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 04 12:17:22 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.594	152	2185	0.400	ng/ul	0.00
4) Naphthalene-d8	10.352	136	8182	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.224	164	4796	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.971	188	9288	0.400	ng/ul	0.00
17) Chrysene-d12	21.164	240	7616	0.400	ng/ul	0.00
23) Perylene-d12	23.321	264	7579	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.189	96	3867	1.594	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	11.948	152	1539	0.112	ng/ul	0.00
18) Fluoranthene-d10	19.001	212	2455	0.097	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.401	128	530	0.022	ng/ul#	72
10) Acenaphthylene	13.941	152	1176	0.058	ng/ul#	87
15) Phenanthrene	17.013	178	1784	0.060	ng/ul#	92
16) Anthracene	17.103	178	1243	0.045	ng/ul#	86
19) Fluoranthene	19.031	202	4101	0.126	ng/ul#	93
21) Benzo(a)anthracene	21.147	228	3286	0.107	ng/ul#	87
22) Chrysene	21.200	228	2238	0.070	ng/ul#	90
24) Benzo(b)fluoranthene	22.682	252	2716	0.080	ng/ul#	50
25) Benzo(k)fluoranthene	22.721	252	1391	0.039	ng/ul#	45

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

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