

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072321\
 Data File : BM031101.D
 Acq On : 23 Jul 2021 21:49
 Operator : CG/JU
 Sample : M2970-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGE81

Manual Integrations
 APPROVED

mohammad
 7/26/2021 1:26:12 PM

Quant Time: Jul 24 00:53:51 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 23 18:04:27 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.639	152	1421	0.40	ng/ul	0.00
4) Naphthalene-d8	10.401	136	5700	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.265	164	3926	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.009	188	8174	0.40	ng/ul	0.00
17) Chrysene-d12	21.195	240	7206	0.40	ng/ul	0.00
23) Perylene-d12	23.370	264	6741	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	7114	4.77	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.993	152	3656	0.37	ng/ul	0.00
18) Fluoranthene-d10	19.039	212	9139	0.42	ng/ul	0.00
Target Compounds						Qvalue
11) Acenaphthene	14.326	153	721	0.05	ng/ul	95
15) Phenanthrene	17.051	178	708	0.03	ng/ul#	80
19) Fluoranthene	19.070	202	1579	0.05	ng/ul#	84
20) Pyrene	19.432	202	1224	0.04	ng/ul#	71
21) Benzo(a)anthracene	21.178	228	764	0.03	ng/ul#	80
22) Chrysene	21.231	228	786	0.03	ng/ul#	83
24) Benzo(b)fluoranthene	22.721	252	975m	0.03	ng/ul	

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

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