

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062821\
 Data File : BM030690.D
 Acq On : 29 Jun 2021 14:22
 Operator : CG/JU
 Sample : M2761-01
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JLX83

Manual Integrations
 APPROVED

mohammad
 6/30/2021 10:25:05 AM

Quant Time: Jun 29 15:00:07 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM062821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 28 14:33:12 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	2814	0.400	ng/ul	0.00
4) Naphthalene-d8	10.571	136	10606	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.413	164	6506	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.151	188	12242m	0.400	ng/ul	0.00
17) Chrysene-d12	21.327	240	10974	0.400	ng/ul	0.00
23) Perylene-d12	23.579	264	13416	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	9894	3.227	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.163	152	6790	0.407	ng/ul	0.00
18) Fluoranthene-d10	19.176	212	15065	0.479	ng/ul	0.00
Target Compounds						
					Qvalue	
10) Acenaphthylene	14.132	152	992	0.034	ng/ul#	72
15) Phenanthrene	17.193	178	2183	0.053	ng/ul#	92
16) Anthracene	17.283	178	796	0.022	ng/ul#	75
19) Fluoranthene	19.206	202	4646	0.096	ng/ul	98
20) Pyrene	19.568	202	3514	0.072	ng/ul#	90
21) Benzo(a)anthracene	21.312	228	1559	0.036	ng/ul#	49
22) Chrysene	21.363	228	1422	0.029	ng/ul#	76
24) Benzo(b)fluoranthene	22.901	252	1377m	0.023	ng/ul	

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

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