

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081221\
 Data File : BM031724.D
 Acq On : 13 Aug 2021 20:51
 Operator : CG/JU
 Sample : M3208-02
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C00E2

Manual Integrations
 APPROVED

mohammad
 8/16/2021 8:40:33 AM

Quant Time: Aug 14 00:51:37 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 : Thu Aug 12 17:28:09 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.566	152	1009	0.400	ng/ul	0.00
4) Naphthalene-d8	10.324	136	3774	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.201	164	2173	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.954	188	4002	0.400	ng/ul	#-0.01
17) Chrysene-d12	21.151	240	3247	0.400	ng/ul	# 0.00
23) Perylene-d12	23.302	264	3492	0.400	ng/ul	#-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.182	96	3608	3.220	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.925	152	1258	0.198	ng/ul	0.00
18) Fluoranthene-d10	18.985	212	2512	0.232	ng/ul	-0.01
Target Compounds						
					Qvalue	
10) Acenaphthylene	13.922	152	317	0.034	ng/ul#	54
15) Phenanthrene	16.995	178	350	0.028	ng/ul#	62
19) Fluoranthene	19.016	202	627	0.045	ng/ul#	60
20) Pyrene	19.378	202	538	0.038	ng/ul#	43
21) Benzo(a)anthracene	21.137	228	367	0.028	ng/ul#	40
22) Chrysene	21.185	228	412	0.030	ng/ul#	56
24) Benzo(b)fluoranthene	22.667	252	485m	0.031	ng/ul	
26) Benzo(a)pyrene	23.210	252	308	0.022	ng/ul#	1

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

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