

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062521\
 Data File : BM030625.D
 Acq On : 26 Jun 2021 06:40
 Operator : CG/JU
 Sample : M2682-04
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGD62

Manual Integrations
 APPROVED

mohammad
 6/28/2021 1:22:15 PM

Quant Time: Jun 26 07:17:10 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM062321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 26 05:08:32 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.794	152	3039	0.400	ng/ul	0.00
4) Naphthalene-d8	10.577	136	12479	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.421	164	7833	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.155	188	15481	0.400	ng/ul	0.00
17) Chrysene-d12	21.323	240	11142	0.400	ng/ul	0.00
23) Perylene-d12	23.576	264	11099	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	6518	2.005	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.173	152	2763	0.138	ng/ul	0.00
18) Fluoranthene-d10	19.180	212	6377	0.187	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	17.197	178	2700	0.049	ng/ul#	91
19) Fluoranthene	19.211	202	5172	0.101	ng/ul	96
20) Pyrene	19.569	202	4796	0.091	ng/ul#	92
21) Benzo(a)anthracene	21.309	228	2531	0.057	ng/ul#	86
22) Chrysene	21.362	228	2839	0.058	ng/ul#	86
24) Benzo(b)fluoranthene	22.897	252	3577m	0.074	ng/ul	
25) Benzo(k)fluoranthene	22.936	252	1355m	0.027	ng/ul	
26) Benzo(a)pyrene	23.474	252	1558	0.036	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	25.858	276	1378	0.025	ng/ul#	97

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

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