

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062521\
 Data File : BM030595.D
 Acq On : 25 Jun 2021 11:05
 Operator : CG/JU
 Sample : M2662-25DL 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGDA1DL

Manual Integrations
 APPROVED

mohammad
 6/28/2021 1:11:18 PM

Quant Time: Jun 25 11:41:30 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 : Fri Jun 25 10:09:59 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.798	152	4215	0.400	ng/ul	0.00
4) Naphthalene-d8	10.586	136	16648	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.425	164	9754	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.159	188	18735	0.400	ng/ul	0.00
17) Chrysene-d12	21.326	240	12589	0.400	ng/ul	0.00
23) Perylene-d12	23.583	264	10754	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.282	96	3067	0.680	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.178	152	1133	0.043	ng/ul	0.00
18) Fluoranthene-d10	19.180	212	2371	0.061	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.635	128	2736	0.055	ng/ul#	92
10) Acenaphthylene	14.144	152	941	0.021	ng/ul#	80
11) Acenaphthene	14.489	153	1014	0.027	ng/ul	97
12) Fluorene	15.471	166	1534	0.037	ng/ul#	97
15) Phenanthrene	17.200	178	13228	0.198	ng/ul	99
16) Anthracene	17.290	178	7088	0.121	ng/ul#	95
19) Fluoranthene	19.210	202	47619	0.823	ng/ul	100
20) Pyrene	19.573	202	29290	0.494	ng/ul	100
21) Benzo(a)anthracene	21.309	228	19157	0.383	ng/ul	96
22) Chrysene	21.360	228	15843	0.285	ng/ul	96
24) Benzo(b)fluoranthene	22.907	252	16702m	0.357	ng/ul	
25) Benzo(k)fluoranthene	22.943	252	6576m	0.134	ng/ul	
26) Benzo(a)pyrene	23.479	252	5110	0.123	ng/ul#	57
27) Indeno(1,2,3-cd)pyrene	25.863	276	3846	0.073	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.870	278	1605	0.038	ng/ul#	17

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

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