

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062521\
 Data File : BM030622.D
 Acq On : 26 Jun 2021 04:51
 Operator : CG/JU
 Sample : M2682-06DL 5X
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGDE4DL

Manual Integrations
 APPROVED

mohammad
 6/28/2021 1:22:07 PM

Quant Time: Jun 26 06:28:46 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	4386	0.400	ng/ul	0.00
4) Naphthalene-d8	10.581	136	18143	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.421	164	11182	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.159	188	22387	0.400	ng/ul	0.00
17) Chrysene-d12	21.321	240	16820	0.400	ng/ul	0.00
23) Perylene-d12	23.573	264	15333	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.283	96	1619	0.345	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.173	152	675	0.023	ng/ul	0.00
18) Fluoranthene-d10	19.180	212	1698	0.033	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.631	128	13183	0.244	ng/ul	99
7) 2-Methylnaphthalene	12.250	142	2056	0.056	ng/ul	100
8) 1-Methylnaphthalene	12.466	142	1252	0.035	ng/ul	100
10) Acenaphthylene	14.140	152	9550	0.182	ng/ul	97
11) Acenaphthene	14.485	153	8124	0.191	ng/ul	98
12) Fluorene	15.467	166	8357	0.177	ng/ul	97
15) Phenanthrene	17.197	178	74720	0.937	ng/ul	100
16) Anthracene	17.291	178	57966	0.830	ng/ul	99
19) Fluoranthene	19.207	202	537474	6.951	ng/ul	98
20) Pyrene	19.569	202	369487	4.664	ng/ul	98
21) Benzo(a)anthracene	21.307	228	269270	4.030	ng/ul	97
22) Chrysene	21.357	228	220841	2.977	ng/ul	98
24) Benzo(b)fluoranthene	22.895	252	244579m	3.667	ng/ul	
25) Benzo(k)fluoranthene	22.939	252	86161m	1.232	ng/ul	
26) Benzo(a)pyrene	23.474	252	114858	1.941	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	25.853	276	72352	0.961	ng/ul#	93
28) Dibenzo(a,h)anthracene	25.858	278	24846	0.414	ng/ul	94
29) Benzo(g,h,i)perylene	26.558	276	3873	0.060	ng/ul#	71

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

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