

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062421\
 Data File : BM030575.D
 Acq On : 24 Jun 2021 14:59
 Operator : CG/JU
 Sample : M2662-23DL2 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGD97DL2

Manual Integrations
 APPROVED

mohammad
 6/25/2021 1:08:22 PM

Quant Time: Jun 24 15:34:02 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 : Thu Jun 24 12:48:41 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.801	152	3934	0.400	ng/ul	0.00
4) Naphthalene-d8	10.590	136	15210	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.428	164	9022	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.162	188	17743	0.400	ng/ul	0.00
17) Chrysene-d12	21.328	240	14025	0.400	ng/ul	0.00
23) Perylene-d12	23.588	264	11772	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.286	96	960	0.228	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.187	152	363	0.015	ng/ul	0.00
18) Fluoranthene-d10	19.184	212	891	0.021	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.640	128	1492	0.033	ng/ul#	85
11) Acenaphthene	14.489	153	1613	0.047	ng/ul	97
12) Fluorene	15.475	166	1745	0.046	ng/ul#	94
15) Phenanthrene	17.207	178	30172	0.478	ng/ul	100
16) Anthracene	17.297	178	7978	0.144	ng/ul	96
19) Fluoranthene	19.214	202	56614	0.878	ng/ul	98
20) Pyrene	19.577	202	35768	0.541	ng/ul	99
21) Benzo(a)anthracene	21.314	228	24206	0.434	ng/ul	97
22) Chrysene	21.365	228	20947	0.339	ng/ul	96
24) Benzo(b)fluoranthene	22.910	252	24031m	0.469	ng/ul	
25) Benzo(k)fluoranthene	22.946	252	8815m	0.164	ng/ul	
26) Benzo(a)pyrene	23.484	252	6765	0.149	ng/ul#	71
27) Indeno(1,2,3-cd)pyrene	25.868	276	5187	0.090	ng/ul#	95
28) Dibenzo(a,h)anthracene	25.873	278	2010	0.044	ng/ul#	35

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

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