

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062521\
 Data File : BM030594.D
 Acq On : 25 Jun 2021 10:29
 Operator : CG/JU
 Sample : M2662-17DL 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGDE6DL

Manual Integrations
 APPROVED

mohammad
 6/28/2021 1:11:08 PM

Quant Time: Jun 25 11:06:23 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	3981	0.400	ng/ul	0.00
4) Naphthalene-d8	10.585	136	15796	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.424	164	9252	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.159	188	18488	0.400	ng/ul	0.00
17) Chrysene-d12	21.326	240	13726	0.400	ng/ul	0.00
23) Perylene-d12	23.583	264	11062	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.282	96	2317	0.544	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.173	152	836	0.033	ng/ul	0.00
18) Fluoranthene-d10	19.180	212	1975	0.047	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.635	128	1816	0.039	ng/ul#	90
15) Phenanthrene	17.200	178	7483	0.114	ng/ul	96
16) Anthracene	17.290	178	4651	0.081	ng/ul#	93
19) Fluoranthene	19.210	202	43543	0.690	ng/ul	100
20) Pyrene	19.573	202	25549	0.395	ng/ul	99
21) Benzo(a)anthracene	21.311	228	23259	0.427	ng/ul	97
22) Chrysene	21.362	228	19858	0.328	ng/ul	97
24) Benzo(b)fluoranthene	22.907	252	20483m	0.426	ng/ul	
25) Benzo(k)fluoranthene	22.946	252	8388m	0.166	ng/ul	
26) Benzo(a)pyrene	23.481	252	6216	0.146	ng/ul#	65
27) Indeno(1,2,3-cd)pyrene	25.863	276	4389	0.081	ng/ul#	95
28) Dibenzo(a,h)anthracene	25.870	278	1865	0.043	ng/ul#	30

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

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