

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072921\
 Data File : BM031360.D
 Acq On : 01 Aug 2021 22:15
 Operator : CG/JU
 Sample : M3106-09DL 5X
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BGD20DL

Manual Integrations
 APPROVED

mohammad
 8/2/2021 3:43:39 PM

Quant Time: Aug 02 01:18:40 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 30 14:16:52 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.604	152	2172	0.400	ng/ul	0.00
4) Naphthalene-d8	10.364	136	8016	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.235	164	4480	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.981	188	8651	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.172	240	6770	0.400	ng/ul	# 0.00
23) Perylene-d12	23.337	264	6510	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	566	0.235	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.961	152	251	0.019	ng/ul	0.00
18) Fluoranthene-d10	19.012	212	547	0.024	ng/ul	0.00
Target Compounds						
					Qvalue	
11) Acenaphthene	14.295	153	756	0.047	ng/ul	99
12) Fluorene	15.289	166	1307	0.071	ng/ul#	98
15) Phenanthrene	17.023	178	31224	1.137	ng/ul	98
16) Anthracene	17.113	178	6557	0.255	ng/ul	99
19) Fluoranthene	19.042	202	54190	1.875	ng/ul#	96
20) Pyrene	19.408	202	40415	1.378	ng/ul	97
21) Benzo(a)anthracene	21.157	228	17905	0.655	ng/ul	94
22) Chrysene	21.208	228	17882m	0.632	ng/ul	
24) Benzo(b)fluoranthene	22.695	252	21832m	0.744	ng/ul	
25) Benzo(k)fluoranthene	22.731	252	7973m	0.263	ng/ul	
26) Benzo(a)pyrene	23.240	252	12500	0.487	ng/ul#	83
27) Indeno(1,2,3-cd)pyrene	25.479	276	9579	0.288	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.486	278	2605	0.097	ng/ul#	17
29) Benzo(g,h,i)perylene	26.135	276	1447	0.051	ng/ul#	24

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

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