

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062923\
 Data File : BM040457.D
 Acq On : 30 Jun 2023 03:16
 Operator : MA/JU
 Sample : 03143-05DL 5X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DCFU4DL

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 06/30/2023
 Supervised By :mohammad ahmed 06/30/2023

Quant Time: Jun 30 04:51:05 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 23 00:44:24 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	7579	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.750	136	27103	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.582	164	12910	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.328	188	24236	0.400	ng/ul	0.00
17) Chrysene-d12	21.509	240	13939	0.400	ng/ul	0.00
23) Perylene-d12	23.920	264	13023	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.329	96	4787	0.402	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.340	152	1099	0.029	ng/ul	-0.01
18) Fluoranthene-d10	19.352	212	1626	0.036	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	17.370	178	6952	0.092	ng/ul	95
16) Anthracene	17.463	178	1720	0.027	ng/ul#	77
19) Fluoranthene	19.384	202	14049	0.237	ng/ul	99
20) Pyrene	19.747	202	11780m	0.186	ng/ul	
21) Benzo(a)anthracene	21.491	228	5640	0.120	ng/ul#	92
22) Chrysene	21.547	228	8439	0.157	ng/ul#	94
24) Benzo(b)fluoranthene	23.186	252	11053	0.219	ng/ul#	68
25) Benzo(k)fluoranthene	23.227	252	3669m	0.074	ng/ul	
26) Benzo(a)pyrene	23.812	252	6098	0.134	ng/ul#	62
27) Indeno(1,2,3-cd)pyrene	26.420	276	5698	0.090	ng/ul#	85
28) Dibenzo(a,h)anthracene	26.430	278	1396	0.028	ng/ul#	1
29) Benzo(g,h,i)perylene	27.198	276	5725	0.103	ng/ul#	83

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

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