

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071423\
 Data File : BM040898.D
 Acq On : 15 Jul 2023 03:06
 Operator : MA/JU
 Sample : 03437-17
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A46H1

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 07/15/2023
 Supervised By :mohammad ahmed 07/17/2023

Quant Time: Jul 15 04:36:48 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 : Sat Jul 15 03:58:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	4542	0.400	ng/ul	0.00
4) Naphthalene-d8	10.658	136	16340	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.504	164	7938	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.257	188	16465	0.400	ng/ul	0.00
17) Chrysene-d12	21.445	240	10748	0.400	ng/ul	0.00
23) Perylene-d12	23.822	264	8843	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	14352	2.012	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.253	152	3426	0.150	ng/ul	0.00
18) Fluoranthene-d10	19.287	212	6879	0.200	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.707	128	18571	0.412	ng/ul	99
7) 2-Methylnaphthalene	12.324	142	973	0.037	ng/ul	98
15) Phenanthrene	17.295	178	7143	0.140	ng/ul	97
16) Anthracene	17.388	178	947	0.022	ng/ul#	74
19) Fluoranthene	19.315	202	15020	0.329	ng/ul	98
20) Pyrene	19.682	202	9362m	0.191	ng/ul	
21) Benzo(a)anthracene	21.431	228	5529	0.152	ng/ul#	92
22) Chrysene	21.483	228	6237	0.150	ng/ul#	92
24) Benzo(b)fluoranthene	23.099	252	7346	0.215	ng/ul#	61
25) Benzo(k)fluoranthene	23.143	252	2849m	0.084	ng/ul	
26) Benzo(a)pyrene	23.713	252	1144	0.037	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	26.277	276	1867	0.043	ng/ul#	85

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

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