

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071124\
 Data File : BM046523.D
 Acq On : 11 Jul 2024 18:29
 Operator : MA/JU
 Sample : P3035-10
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4CX8

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 07/12/2024
 Supervised By :mohammad ahmed 07/13/2024

Quant Time: Jul 12 04:22:44 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM071024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 04:18:37 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.514	152	2770	0.400	ng/ul	0.00
4) Naphthalene-d8	10.276	136	7432	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.159	164	5016	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.908	188	11080	0.400	ng/ul	0.00
17) Chrysene-d12	21.114	240	9641	0.400	ng/ul	0.00
23) Perylene-d12	23.259	264	10721m	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.122	96	4599	2.021	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.876	152	2026	0.169	ng/ul	0.00
18) Fluoranthene-d10	18.945	212	5451	0.186	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	16.950	178	6432	0.203	ng/ul	99
16) Anthracene	17.043	178	1342	0.043	ng/ul#	87
19) Fluoranthene	18.978	202	14846	0.375	ng/ul	99
20) Pyrene	19.340	202	14303	0.334	ng/ul	99
21) Benzo(a)anthracene	21.097	228	7346	0.172	ng/ul	96
22) Chrysene	21.149	228	7852	0.190	ng/ul	98
24) Benzo(b)fluoranthene	22.625	252	11370m	0.266	ng/ul	
25) Benzo(k)fluoranthene	22.660	252	4217m	0.100	ng/ul	
26) Benzo(a)pyrene	23.169	252	7645	0.217	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	25.380	276	6754	0.126	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.401	278	1632	0.038	ng/ul#	23
29) Benzo(g,h,i)perylene	26.028	276	6801	0.159	ng/ul#	94

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

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