

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071924\
 Data File : BM046735.D
 Acq On : 19 Jul 2024 20:29
 Operator : MA/JU
 Sample : P3201-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4BZ2

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/20/2024
 Supervised By :mohammad ahmed 07/20/2024

Quant Time: Jul 19 23:58:41 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 19 11:38:04 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.484	152	4032	0.400	ng/ul	0.00
4) Naphthalene-d8	10.248	136	11269	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.131	164	6667	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.887	188	12829m	0.400	ng/ul	0.00
17) Chrysene-d12	21.093	240	9190	0.400	ng/ul #	0.00
23) Perylene-d12	23.227	264	9881	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.105	96	8837	2.668	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.842	152	3870	0.213	ng/ul	0.00
18) Fluoranthene-d10	18.926	212	6355	0.227	ng/ul	0.00
Target Compounds						
						Qvalue
5) Naphthalene	10.292	128	2480	0.087	ng/ul	95
7) 2-Methylnaphthalene	11.919	142	910	0.045	ng/ul	99
8) 1-Methylnaphthalene	12.139	142	700	0.034	ng/ul	99
10) Acenaphthylene	13.848	152	2733	0.090	ng/ul	95
11) Acenaphthene	14.195	153	1892	0.093	ng/ul	98
12) Fluorene	15.190	166	2463	0.098	ng/ul	93
15) Phenanthrene	16.929	178	37817	1.033	ng/ul	99
16) Anthracene	17.017	178	7806	0.218	ng/ul	99
19) Fluoranthene	18.954	202	66685	1.767	ng/ul	99
20) Pyrene	19.321	202	57395	1.404	ng/ul	99
21) Benzo(a)anthracene	21.079	228	28159	0.693	ng/ul	96
22) Chrysene	21.129	228	29808	0.758	ng/ul	98
24) Benzo(b)fluoranthene	22.598	252	42631m	1.080	ng/ul	
25) Benzo(k)fluoranthene	22.633	252	12262m	0.316	ng/ul	
26) Benzo(a)pyrene	23.136	252	26161	0.805	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	25.340	276	21992	0.445	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.350	278	5231	0.134	ng/ul#	62
29) Benzo(g,h,i)perylene	25.984	276	22252	0.564	ng/ul	100

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

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