

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM050324\
 Data File : BM045697.D
 Acq On : 03 May 2024 15:51
 Operator : MA/JU
 Sample : P2238-10MS
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E1CX6MS

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 05/03/2024
 Supervised By :mohammad ahmed 05/06/2024

Quant Time: May 03 16:56:32 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM050324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 03 03:27:04 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.551	152	237	0.400	ng/ul	-0.03
4) Naphthalene-d8	10.298	136	812	0.400	ng/ul	#-0.05
9) Acenaphthene-d10	14.177	164	632	0.400	ng/ul	-0.02
13) Phenanthrene-d10	16.942	188	1406	0.400	ng/ul	#-0.03
17) Chrysene-d12	21.158	240	1765	0.400	ng/ul	#-0.03
23) Perylene-d12	23.332	264	2488	0.400	ng/ul	-0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.150	96	129	0.481	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.898	152	295	0.230	ng/ul	-0.05
18) Fluoranthene-d10	18.982	212	976	0.220	ng/ul	-0.03
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.180	88	252	0.906	ng/ul	89
5) Naphthalene	10.347	128	1217	0.588	ng/ul#	79
7) 2-Methylnaphthalene	11.975	142	692	0.553	ng/ul	98
8) 1-Methylnaphthalene	12.189	142	532	0.384	ng/ul	98
10) Acenaphthylene	13.900	152	875	0.330	ng/ul	95
11) Acenaphthene	14.242	153	1684	0.855	ng/ul	96
12) Fluorene	15.255	166	1317	0.577	ng/ul#	92
14) Pentachlorophenol	16.625	266	11	0.044	ng/ul#	58
15) Phenanthrene	16.984	178	13912m	3.984	ng/ul	
16) Anthracene	17.085	178	3207	0.946	ng/ul#	89
19) Fluoranthene	19.010	202	25089	4.473	ng/ul#	84
20) Pyrene	19.378	202	23121	3.876	ng/ul#	85
21) Benzo(a)anthracene	21.144	228	19838	3.851	ng/ul#	91
22) Chrysene	21.193	228	23610	3.238	ng/ul	93
24) Benzo(b)fluoranthene	22.683	252	61013m	8.804	ng/ul	
25) Benzo(k)fluoranthene	22.724	252	22418m	2.613	ng/ul	
26) Benzo(a)pyrene	23.236	252	48611	6.329	ng/ul#	74
27) Indeno(1,2,3-cd)pyrene	25.491	276	49186	4.649	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.504	278	12654m	1.496	ng/ul	
29) Benzo(g,h,i)perylene	26.152	276	55286	5.894	ng/ul#	83

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

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