

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100423\
 Data File : BM042196.D
 Acq On : 04 Oct 2023 21:14
 Operator : MA/JU
 Sample : 04588-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BBHT7

Manual Integrations
 APPROVED

Reviewed By :Yogesh
 Patel

Quant Time: Oct 05 01:55:17 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM100223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 02 19:28:41 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							10/05/2023
1) 1,4-Dichlorobenzene-d4	7.950	152	4023	0.400	ng/ul	# 0.01	Supervised By :mohammad ahmed
4) Naphthalene-d8	10.746	136	7491	0.400	ng/ul	# 0.00	
9) Acenaphthene-d10	14.578	164	5454	0.400	ng/ul	# 0.00	
13) Phenanthrene-d10	17.329	188	8053	0.400	ng/ul	#-0.01	
17) Chrysene-d12	21.504	240	4999	0.400	ng/ul	-0.01	
23) Perylene-d12	23.901	264	5223	0.400	ng/ul	-0.02	10/06/2023
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.360	96	14387	2.685	ng/ul	0.00	
6) 2-Methylnaphthalene-d10	12.324	152	1929	0.188	ng/ul	-0.01	
18) Fluoranthene-d10	19.352	212	1541	0.082	ng/ul	0.00	
Target Compounds							Qvalue
2) 1,4-Dioxane	3.398	88	825	0.147	ng/ul#	69	
5) Naphthalene	10.796	128	181963	7.281	ng/ul#	88	
7) 2-Methylnaphthalene	12.401	142	426088	28.475	ng/ul	98	
8) 1-Methylnaphthalene	12.621	142	413020	27.402	ng/ul	100	
10) Acenaphthylene	14.291	152	8649	0.251	ng/ul#	1	
11) Acenaphthene	14.643	153	23520	0.902	ng/ul#	77	
12) Fluorene	15.628	166	42753	1.455	ng/ul#	91	
15) Phenanthrene	17.371	178	69750	2.101	ng/ul#	95	
16) Anthracene	17.464	178	4514	0.149	ng/ul#	64	
19) Fluoranthene	19.380	202	5824	0.153	ng/ul#	83	
20) Pyrene	19.747	202	12487m	0.309	ng/ul		
21) Benzo(a)anthracene	21.486	228	4124	0.128	ng/ul#	82	
22) Chrysene	21.539	228	5060m	0.161	ng/ul		
24) Benzo(b)fluoranthene	23.164	252	3446	0.112	ng/ul#	39	
25) Benzo(k)fluoranthene	23.208	252	1471m	0.049	ng/ul		
26) Benzo(a)pyrene	23.792	252	2506	0.094	ng/ul#	39	
27) Indeno(1,2,3-cd)pyrene	26.391	276	1844	0.049	ng/ul#	95	
28) Dibenzo(a,h)anthracene	26.394	278	1252	0.044	ng/ul#	4	
29) Benzo(g,h,i)perylene	27.156	276	2913	0.093	ng/ul#	79	

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

Manual Integrations

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