

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012623\
 Data File : BM038585.D
 Acq On : 27 Jan 2023 14:25
 Operator : CG/JU
 Sample : 01226-10MSD
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EX9G4MSD

Quant Time: Jan 27 21:59:54 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM012623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 27 00:18:15 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.962	152	1291	0.400	ng/ul	0.00
4) Naphthalene-d8	10.774	136	3769	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.597	164	2248	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.341	188	4903	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.512	240	5260	0.400	ng/ul	# 0.00
23) Perylene-d12	23.921	264	5894	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.343	96	245	0.145	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.357	152	1337	0.229	ng/ul	0.00
18) Fluoranthene-d10	19.362	212	4143	0.276	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.376	88	982	0.571	ng/ul#	25
5) Naphthalene	10.823	128	2523	0.206	ng/ul	98
7) 2-Methylnaphthalene	12.429	142	1620	0.250	ng/ul	95
8) 1-Methylnaphthalene	12.649	142	1723	0.258	ng/ul	100
10) Acenaphthylene	14.319	152	2847	0.287	ng/ul	93
11) Acenaphthene	14.657	153	2011	0.266	ng/ul	98
12) Fluorene	15.642	166	2491	0.264	ng/ul#	96
14) Pentachlorophenol	16.991	266	1182	0.592	ng/ul	99
15) Phenanthrene	17.379	178	4331	0.273	ng/ul#	94
16) Anthracene	17.472	178	4509	0.334	ng/ul	94
19) Fluoranthene	19.389	202	5815	0.310	ng/ul#	91
20) Pyrene	19.757	202	6311	0.313	ng/ul#	96
21) Benzo(a)anthracene	21.498	228	5877	0.314	ng/ul	93
22) Chrysene	21.550	228	5813	0.310	ng/ul	93
24) Benzo(b)fluoranthene	23.184	252	6643	0.299	ng/ul	88
25) Benzo(k)fluoranthene	23.234	252	6446	0.299	ng/ul#	88
26) Benzo(a)pyrene	23.816	252	5685	0.285	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	26.428	276	8084	0.288	ng/ul#	65
28) Dibenzo(a,h)anthracene	26.444	278	6577	0.295	ng/ul#	91
29) Benzo(g,h,i)perylene	27.202	276	6936	0.285	ng/ul	97

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

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