

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070823\
 Data File : BN026396.D
 Acq On : 08 Jul 2023 21:40
 Operator : MA/JU
 Sample : 03350-03MS
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 YBW27MS

Quant Time: Jul 08 23:54:14 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 08 09:20:18 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.930	152	5210	0.400	ng/ul	0.00
4) Naphthalene-d8	10.737	136	18433	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.570	164	9590	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.322	188	17419	0.400	ng/ul	0.00
17) Chrysene-d12	21.514	240	10859	0.400	ng/ul	-0.01
23) Perylene-d12	23.961	264	10553	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	2683	0.457	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.327	152	6337	0.224	ng/ul	0.00
18) Fluoranthene-d10	19.354	212	10595	0.256	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.335	88	6741	1.082	ng/ul#	87
5) Naphthalene	10.787	128	13481	0.255	ng/ul	99
7) 2-Methylnaphthalene	12.404	142	7827	0.233	ng/ul	100
8) 1-Methylnaphthalene	12.618	142	8412	0.241	ng/ul	100
10) Acenaphthylene	14.293	152	10311	0.230	ng/ul	98
11) Acenaphthene	14.630	153	10016	0.273	ng/ul	98
12) Fluorene	15.620	166	10323	0.245	ng/ul	98
14) Pentachlorophenol	16.984	266	1177	0.189	ng/ul	99
15) Phenanthrene	17.364	178	14312	0.249	ng/ul	98
16) Anthracene	17.457	178	13925	0.272	ng/ul	96
19) Fluoranthene	19.382	202	14990	0.258	ng/ul	100
20) Pyrene	19.745	202	16028	0.266	ng/ul	98
21) Benzo(a)anthracene	21.500	228	10877	0.243	ng/ul	99
22) Chrysene	21.552	228	12059	0.259	ng/ul	99
24) Benzo(b)fluoranthene	23.215	252	11609	0.258	ng/ul	86
25) Benzo(k)fluoranthene	23.262	252	11706	0.259	ng/ul#	88
26) Benzo(a)pyrene	23.862	252	9968	0.255	ng/ul#	81
27) Indeno(1,2,3-cd)pyrene	26.527	276	12579	0.294	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.534	278	9596	0.297	ng/ul	92
29) Benzo(g,h,i)perylene	27.312	276	11261	0.296	ng/ul	98

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

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