

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062922\
 Data File : BN020575.D
 Acq On : 29 Jun 2022 18:31
 Operator : CG/JU
 Sample : N3355-12MS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EY0C9MS

Quant Time: Jun 29 19:06:12 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 24 00:17:34 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.741	152	2868	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.524	136	9631	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.368	164	5518	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.112	188	11202	0.400	ng/ul	0.00
17) Chrysene-d12	21.303	240	9011	0.400	ng/ul	0.00
23) Perylene-d12	23.586	264	6678	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	964	0.290	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.119	152	5412	0.342	ng/ul	0.00
18) Fluoranthene-d10	19.141	212	13224	0.445	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.227	88	2694	0.794	ng/ul#	81
5) Naphthalene	10.568	128	10195	0.343	ng/ul	99
7) 2-Methylnaphthalene	12.190	142	6375	0.320	ng/ul	99
8) 1-Methylnaphthalene	12.410	142	6805	0.361	ng/ul	100
10) Acenaphthylene	14.086	152	9383	0.359	ng/ul	99
11) Acenaphthene	14.433	153	7355	0.351	ng/ul	98
12) Fluorene	15.418	166	8488	0.342	ng/ul	99
14) Pentachlorophenol	16.778	266	1444	0.639	ng/ul	100
15) Phenanthrene	17.154	178	15204	0.390	ng/ul	99
16) Anthracene	17.247	178	11800	0.351	ng/ul	100
19) Fluoranthene	19.174	202	17299	0.420	ng/ul	99
20) Pyrene	19.537	202	17593	0.412	ng/ul	99
21) Benzo(a)anthracene	21.286	228	12926	0.399	ng/ul	100
22) Chrysene	21.341	228	15253	0.423	ng/ul	100
24) Benzo(b)fluoranthene	22.893	252	12971	0.420	ng/ul	98
25) Benzo(k)fluoranthene	22.940	252	12928	0.414	ng/ul	99
26) Benzo(a)pyrene	23.484	252	10395	0.370	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.916	276	11295	0.404	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.926	278	9228	0.409	ng/ul	100
29) Benzo(g,h,i)perylene	26.627	276	9892	0.412	ng/ul	100

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

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