

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071122\
 Data File : BN020703.D
 Acq On : 11 Jul 2022 15:25
 Operator : CG/JU
 Sample : N3577-03MS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 YB324MS

Quant Time: Jul 11 18:38:25 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 : Sat Jul 02 00:47:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.825	152	4013	0.400	ng/ul	0.00
4) Naphthalene-d8	10.623	136	12747	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.474	164	7010	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.225	188	13383	0.400	ng/ul	0.00
17) Chrysene-d12	21.432	240	9728	0.400	ng/ul	-0.01
23) Perylene-d12	23.785	264	6724	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.227	96	1061	0.228	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.223	152	6106	0.292	ng/ul	0.00
18) Fluoranthene-d10	19.262	212	14287	0.445	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.260	88	3123	0.658	ng/ul	88
5) Naphthalene	10.672	128	12167	0.309	ng/ul	99
7) 2-Methylnaphthalene	12.294	142	7259	0.276	ng/ul	98
8) 1-Methylnaphthalene	12.514	142	7701	0.308	ng/ul	100
10) Acenaphthylene	14.192	152	9998	0.301	ng/ul	100
11) Acenaphthene	14.534	153	8102	0.305	ng/ul	99
12) Fluorene	15.529	166	8977	0.285	ng/ul	98
14) Pentachlorophenol	16.892	266	1397	0.518	ng/ul	99
15) Phenanthrene	17.268	178	15689	0.337	ng/ul	100
16) Anthracene	17.361	178	12545	0.312	ng/ul	100
19) Fluoranthene	19.295	202	18962	0.427	ng/ul	100
20) Pyrene	19.657	202	19431	0.422	ng/ul	98
21) Benzo(a)anthracene	21.417	228	12467	0.357	ng/ul	100
22) Chrysene	21.470	228	15089	0.388	ng/ul	100
24) Benzo(b)fluoranthene	23.068	252	12238	0.393	ng/ul	99
25) Benzo(k)fluoranthene	23.118	252	12674	0.403	ng/ul	99
26) Benzo(a)pyrene	23.680	252	10806	0.382	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.218	276	11043	0.392	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.238	278	8887	0.391	ng/ul	98
29) Benzo(g,h,i)perylene	26.959	276	10029	0.415	ng/ul	100

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

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