

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080322\
 Data File : BN020977.D
 Acq On : 03 Aug 2022 14:28
 Operator : CG/JU
 Sample : SSTD0.831
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.8233

Quant Time: Aug 03 14:55:33 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 03 14:49:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	3050	0.400	ng/ul	0.00
4) Naphthalene-d8	10.602	136	8054	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.457	164	3903	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.210	188	7436	0.400	ng/ul	0.00
17) Chrysene-d12	21.421	240	8819	0.400	ng/ul	0.00
23) Perylene-d12	23.766	264	7142	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	3108	0.836	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.202	152	9116	0.768	ng/ul	0.00
18) Fluoranthene-d10	19.249	212	23110	0.856	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.245	88	3251	0.812	ng/ul	90
5) Naphthalene	10.651	128	17960	0.816	ng/ul	98
7) 2-Methylnaphthalene	12.274	142	10110	0.754	ng/ul	99
8) 1-Methylnaphthalene	12.494	142	11059	0.747	ng/ul	100
10) Acenaphthylene	14.175	152	13774	0.863	ng/ul	99
11) Acenaphthene	14.517	153	11167	0.838	ng/ul	99
12) Fluorene	15.512	166	12365	0.805	ng/ul	99
14) Pentachlorophenol	16.876	266	898	0.756	ng/ul	99
15) Phenanthrene	17.248	178	19619	0.778	ng/ul	99
16) Anthracene	17.345	178	16311	0.753	ng/ul	98
19) Fluoranthene	19.277	202	30853	0.778	ng/ul	98
20) Pyrene	19.644	202	30587	0.737	ng/ul	98
21) Benzo(a)anthracene	21.407	228	22199	0.693	ng/ul	100
22) Chrysene	21.456	228	28595	0.722	ng/ul	100
24) Benzo(b)fluoranthene	23.052	252	25960	0.761	ng/ul	97
25) Benzo(k)fluoranthene	23.099	252	27916	0.755	ng/ul	96
26) Benzo(a)pyrene	23.660	252	22905	0.712	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	26.186	276	26520	0.701	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.199	278	21303	0.688	ng/ul	96
29) Benzo(g,h,i)perylene	26.924	276	23243	0.675	ng/ul	99

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

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