

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080922\
 Data File : BN021082.D
 Acq On : 09 Aug 2022 13:03
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4276

Quant Time: Aug 09 23:11:32 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 : Tue Aug 09 23:10:06 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.780	152	3009	0.400	ng/ul	0.00
4) Naphthalene-d8	10.579	136	9747	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.437	164	5715	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.192	188	11192	0.400	ng/ul	0.00
17) Chrysene-d12	21.404	240	8598	0.400	ng/ul	0.00
23) Perylene-d12	23.740	264	6572	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.194	96	1471	0.382	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.179	152	6101	0.427	ng/ul	0.00
18) Fluoranthene-d10	19.230	212	12020	0.453	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.228	88	1564	0.394	ng/ul#	90
5) Naphthalene	10.628	128	10362	0.346	ng/ul	98
7) 2-Methylnaphthalene	12.250	142	6702	0.369	ng/ul	100
8) 1-Methylnaphthalene	12.470	142	7180	0.395	ng/ul	100
10) Acenaphthylene	14.155	152	9556	0.358	ng/ul	99
11) Acenaphthene	14.502	153	7380	0.334	ng/ul	99
12) Fluorene	15.492	166	8264	0.342	ng/ul	99
14) Pentachlorophenol	16.854	266	810	0.469	ng/ul	99
15) Phenanthrene	17.230	178	13020	0.325	ng/ul	99
16) Anthracene	17.322	178	11377	0.367	ng/ul	98
19) Fluoranthene	19.262	202	14269	0.375	ng/ul	99
20) Pyrene	19.625	202	14574	0.371	ng/ul	98
21) Benzo(a)anthracene	21.387	228	10615	0.378	ng/ul	99
22) Chrysene	21.442	228	11601	0.314	ng/ul	98
24) Benzo(b)fluoranthene	23.029	252	10097	0.332	ng/ul	97
25) Benzo(k)fluoranthene	23.076	252	10000	0.311	ng/ul	98
26) Benzo(a)pyrene	23.634	252	9188	0.331	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.149	276	8766	0.265	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.166	278	6952	0.265	ng/ul	98
29) Benzo(g,h,i)perylene	26.887	276	7904	0.267	ng/ul	97

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

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