

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080122\
 Data File : BN020929.D
 Acq On : 01 Aug 2022 20:50
 Operator : CG/JU
 Sample : SST0.825
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.8226

Quant Time: Aug 02 03:36:25 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 02 03:35:30 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	2290	0.400	ng/ul	0.00
4) Naphthalene-d8	10.602	136	7138	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.461	164	4057	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.210	188	7895	0.400	ng/ul	0.00
17) Chrysene-d12	21.418	240	7109	0.400	ng/ul	0.00
23) Perylene-d12	23.768	264	6096	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	2383	0.866	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.202	152	9252	0.853	ng/ul	0.00
18) Fluoranthene-d10	19.249	212	18687	0.823	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.244	88	2569	0.909	ng/ul	90
5) Naphthalene	10.651	128	16342	0.730	ng/ul	98
7) 2-Methylnaphthalene	12.279	142	10280	0.727	ng/ul	99
8) 1-Methylnaphthalene	12.494	142	11184	0.817	ng/ul	100
10) Acenaphthylene	14.179	152	14101	0.699	ng/ul	99
11) Acenaphthene	14.521	153	11602	0.711	ng/ul	98
12) Fluorene	15.511	166	13098	0.731	ng/ul	99
14) Pentachlorophenol	16.876	266	1050	0.883	ng/ul	98
15) Phenanthrene	17.252	178	20671	0.714	ng/ul	99
16) Anthracene	17.349	178	17138	0.739	ng/ul	99
19) Fluoranthene	19.282	202	22970	0.690	ng/ul	97
20) Pyrene	19.644	202	23106	0.665	ng/ul	99
21) Benzo(a)anthracene	21.404	228	17739	0.755	ng/ul	100
22) Chrysene	21.456	228	21491	0.711	ng/ul	99
24) Benzo(b)fluoranthene	23.052	252	20402	0.706	ng/ul	96
25) Benzo(k)fluoranthene	23.099	252	21868	0.732	ng/ul	96
26) Benzo(a)pyrene	23.660	252	19025	0.714	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.189	276	22876	0.816	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.202	278	18416	0.862	ng/ul	95
29) Benzo(g,h,i)perylene	26.927	276	20061	0.794	ng/ul	99

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

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