

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082222\
 Data File : BN021301.D
 Acq On : 22 Aug 2022 14:26
 Operator : CG/JU
 Sample : SST0.235
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.2238

Quant Time: Aug 22 17:53:04 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN082222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 22 17:51:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.093	152	2909	0.400	ng/u1	0.00
4) Naphthalene-d8	10.927	136	9084	0.400	ng/u1	# 0.00
9) Acenaphthene-d10	14.749	164	4684	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.498	188	9544	0.400	ng/u1	# 0.00
17) Chrysene-d12	21.697	240	7666	0.400	ng/u1	# 0.00
23) Perylene-d12	24.231	264	7178	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.385	96	731	0.196	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.511	152	2563	0.193	ng/u1	0.00
18) Fluoranthene-d10	19.529	212	4424	0.187	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.422	88	713	0.186	ng/u1	95
5) Naphthalene	10.977	128	6363	0.228	ng/u1	99
7) 2-Methylnaphthalene	12.588	142	3917	0.231	ng/u1	98
8) 1-Methylnaphthalene	12.803	142	3150	0.186	ng/u1	100
10) Acenaphthylene	14.472	152	5305	0.242	ng/u1	100
11) Acenaphthene	14.814	153	4052	0.224	ng/u1	98
12) Fluorene	15.799	166	4427	0.224	ng/u1	100
14) Pentachlorophenol	17.143	266	222	0.151	ng/u1	96
15) Phenanthrene	17.540	178	7636	0.224	ng/u1	100
16) Anthracene	17.633	178	6017	0.228	ng/u1	99
19) Fluoranthene	19.557	202	7711	0.227	ng/u1	96
20) Pyrene	19.920	202	8061	0.230	ng/u1#	94
21) Benzo(a)anthracene	21.679	228	5699	0.228	ng/u1	100
22) Chrysene	21.732	228	7483	0.227	ng/u1	100
24) Benzo(b)fluoranthene	23.451	252	6096	0.183	ng/u1	96
25) Benzo(k)fluoranthene	23.503	252	6928	0.198	ng/u1#	94
26) Benzo(a)pyrene	24.120	252	6015	0.198	ng/u1#	95
27) Indeno(1,2,3-cd)pyrene	26.911	276	6951	0.192	ng/u1#	94
28) Dibenzo(a,h)anthracene	26.938	278	5755	0.201	ng/u1	94
29) Benzo(g,h,i)perylene	27.733	276	7021	0.217	ng/u1	96

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

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