

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN030923\
 Data File : BN024165.D
 Acq On : 08 Mar 2023 13:08
 Operator : CG/JU
 Sample : SST0.277
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.2238

Quant Time: Mar 09 00:49:55 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN030923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 09 00:47:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.055	152	8725	0.400	ng/ul	0.00
4) Naphthalene-d8	10.878	136	24607	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.694	164	12720	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.434	188	26371	0.400	ng/ul	0.00
17) Chrysene-d12	21.627	240	19093	0.400	ng/ul	0.00
23) Perylene-d12	24.129	264	16183	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.410	96	1914	0.205	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.456	152	5755	0.149	ng/ul	0.00
18) Fluoranthene-d10	19.460	212	10271	0.200	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.448	88	2104	0.212	ng/ul#	90
5) Naphthalene	10.927	128	12239	0.182	ng/ul	99
7) 2-Methylnaphthalene	12.528	142	7325	0.167	ng/ul	99
8) 1-Methylnaphthalene	12.748	142	7736	0.173	ng/ul	100
10) Acenaphthylene	14.411	152	10308	0.212	ng/ul	99
11) Acenaphthene	14.754	153	7998	0.187	ng/ul	99
12) Fluorene	15.739	166	8807	0.170	ng/ul	98
14) Pentachlorophenol	17.080	266	1255	0.144	ng/ul#	73
15) Phenanthrene	17.477	178	14710	0.187	ng/ul	99
16) Anthracene	17.570	178	12595	0.191	ng/ul	99
19) Fluoranthene	19.487	202	15105	0.238	ng/ul	96
20) Pyrene	19.850	202	15478	0.235	ng/ul	96
21) Benzo(a)anthracene	21.609	228	12405	0.191	ng/ul	100
22) Chrysene	21.662	228	13098	0.192	ng/ul	100
24) Benzo(b)fluoranthene	23.363	252	11675	0.163	ng/ul	93
25) Benzo(k)fluoranthene	23.410	252	12485	0.179	ng/ul#	93
26) Benzo(a)pyrene	24.015	252	10228	0.172	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	26.747	276	11376	0.136	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.770	278	8617	0.129	ng/ul	95
29) Benzo(g,h,i)perylene	27.552	276	9928	0.144	ng/ul	97

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

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