

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112622\
 Data File : BN022875.D
 Acq On : 26 Nov 2022 10:33
 Operator : CG/JU
 Sample : SST0.253
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.2210

Quant Time: Nov 28 03:01:01 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN112622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 28 02:35:52 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.069	152	7933	0.400	ng/ul	0.00
4) Naphthalene-d8	10.892	136	21255	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.710	164	11500	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.453	188	24775	0.400	ng/ul	0.00
17) Chrysene-d12	21.638	240	15327	0.400	ng/ul	0.00
23) Perylene-d12	24.126	264	10623	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.412	96	1910	0.201	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.476	152	6447	0.199	ng/ul	0.00
18) Fluoranthene-d10	19.476	212	11910	0.210	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.450	88	1983	0.200	ng/ul#	95
5) Naphthalene	10.942	128	12741	0.198	ng/ul	99
7) 2-Methylnaphthalene	12.547	142	7958	0.197	ng/ul	99
8) 1-Methylnaphthalene	12.767	142	8147	0.199	ng/ul	100
10) Acenaphthylene	14.432	152	11063	0.195	ng/ul	99
11) Acenaphthene	14.770	153	8905	0.198	ng/ul	97
12) Fluorene	15.755	166	10043	0.198	ng/ul	99
14) Pentachlorophenol	17.099	266	886	0.171	ng/ul#	100
15) Phenanthrene	17.496	178	16843	0.198	ng/ul	100
16) Anthracene	17.588	178	13983	0.192	ng/ul	99
19) Fluoranthene	19.508	202	16565	0.209	ng/ul	99
20) Pyrene	19.871	202	16543	0.213	ng/ul	99
21) Benzo(a)anthracene	21.620	228	11357	0.194	ng/ul	100
22) Chrysene	21.673	228	12898	0.201	ng/ul	99
24) Benzo(b)fluoranthene	23.363	252	10289	0.191	ng/ul	94
25) Benzo(k)fluoranthene	23.412	252	10945	0.192	ng/ul#	94
26) Benzo(a)pyrene	24.014	252	7817	0.191	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	26.736	276	9029	0.185	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.766	278	7109	0.184	ng/ul	93
29) Benzo(g,h,i)perylene	27.534	276	8083	0.190	ng/ul	97

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

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