

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080322\
 Data File : BN020976.D
 Acq On : 03 Aug 2022 13:51
 Operator : CG/JU
 Sample : SST0.430
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.4232

Quant Time: Aug 03 14:50:53 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 : Wed Aug 03 14:49:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.814	152	2386	0.400	ng/ul	0.00
4) Naphthalene-d8	10.607	136	8079	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.456	164	5412	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.209	188	10657	0.400	ng/ul	0.00
17) Chrysene-d12	21.418	240	8916	0.400	ng/ul	0.00
23) Perylene-d12	23.765	264	5906	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	1488	0.511	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.207	152	5625	0.472	ng/ul	0.00
18) Fluoranthene-d10	19.249	212	11176	0.410	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.244	88	1519	0.485	ng/ul	95
5) Naphthalene	10.651	128	8984	0.407	ng/ul	99
7) 2-Methylnaphthalene	12.279	142	6447	0.479	ng/ul	98
8) 1-Methylnaphthalene	12.499	142	7167	0.483	ng/ul	100
10) Acenaphthylene	14.179	152	8560	0.387	ng/ul	99
11) Acenaphthene	14.521	153	7320	0.396	ng/ul	100
12) Fluorene	15.516	166	8049	0.378	ng/ul	99
14) Pentachlorophenol	16.880	266	494	0.290	ng/ul	98
15) Phenanthrene	17.252	178	13117	0.363	ng/ul	100
16) Anthracene	17.349	178	10097	0.325	ng/ul	99
19) Fluoranthene	19.282	202	14075	0.351	ng/ul	100
20) Pyrene	19.644	202	14359	0.342	ng/ul	97
21) Benzo(a)anthracene	21.406	228	9970	0.308	ng/ul	99
22) Chrysene	21.459	228	13467	0.336	ng/ul	99
24) Benzo(b)fluoranthene	23.052	252	9017	0.320	ng/ul	99
25) Benzo(k)fluoranthene	23.099	252	9365	0.306	ng/ul	99
26) Benzo(a)pyrene	23.660	252	8215	0.309	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	26.185	276	12321	0.394	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.209	278	9695	0.379	ng/ul	98
29) Benzo(g,h,i)perylene	26.930	276	11207	0.394	ng/ul	98

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

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