

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080222\
 Data File : BN020947.D
 Acq On : 02 Aug 2022 17:37
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4262

Quant Time: Aug 02 22:34:06 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 04:19:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	3029	0.400	ng/ul	0.00
4) Naphthalene-d8	10.607	136	8402	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.461	164	6186	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.214	188	12335	0.400	ng/ul	0.00
17) Chrysene-d12	21.424	240	10491	0.400	ng/ul	0.00
23) Perylene-d12	23.771	264	8575	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.215	96	1645	0.445	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.208	152	6380	0.515	ng/ul	0.00
18) Fluoranthene-d10	19.254	212	12988	0.405	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.249	88	1736	0.437	ng/ul	91
5) Naphthalene	10.651	128	9027	0.393	ng/ul	99
7) 2-Methylnaphthalene	12.279	142	7166	0.512	ng/ul	99
8) 1-Methylnaphthalene	12.499	142	7956	0.515	ng/ul	100
10) Acenaphthylene	14.179	152	9946	0.393	ng/ul	99
11) Acenaphthene	14.521	153	8277	0.392	ng/ul	100
12) Fluorene	15.516	166	9116	0.375	ng/ul	99
14) Pentachlorophenol	16.880	266	657	0.334	ng/ul	99
15) Phenanthrene	17.256	178	14844	0.355	ng/ul	100
16) Anthracene	17.353	178	11730	0.326	ng/ul	99
19) Fluoranthene	19.282	202	16121	0.342	ng/ul	98
20) Pyrene	19.644	202	16549	0.335	ng/ul	94
21) Benzo(a)anthracene	21.410	228	11956	0.314	ng/ul	100
22) Chrysene	21.459	228	15225	0.323	ng/ul	100
24) Benzo(b)fluoranthene	23.058	252	14029	0.343	ng/ul	99
25) Benzo(k)fluoranthene	23.105	252	14773	0.333	ng/ul	99
26) Benzo(a)pyrene	23.666	252	12551	0.325	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.192	276	13918	0.307	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.212	278	11059	0.297	ng/ul	97
29) Benzo(g,h,i)perylene	26.937	276	10719	0.259	ng/ul	99

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

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