

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080422\
 Data File : BN020982.D
 Acq On : 04 Aug 2022 09:53
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4266

Quant Time: Aug 04 11:02:33 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 16:31:11 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	4041	0.400	ng/ul	0.00
4) Naphthalene-d8	10.601	136	9616	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.455	164	6973	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.208	188	13641	0.400	ng/ul	0.00
17) Chrysene-d12	21.416	240	11125	0.400	ng/ul	0.00
23) Perylene-d12	23.757	264	8855	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	1905	0.368	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.201	152	5513	0.391	ng/ul	0.00
18) Fluoranthene-d10	19.248	212	14160	0.412	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.245	88	1861	0.349	ng/ul	92
5) Naphthalene	10.650	128	10273	0.348	ng/ul	99
7) 2-Methylnaphthalene	12.278	142	6124	0.342	ng/ul	100
8) 1-Methylnaphthalene	12.492	142	6831	0.381	ng/ul	100
10) Acenaphthylene	14.173	152	11441	0.351	ng/ul	98
11) Acenaphthene	14.516	153	9512	0.353	ng/ul	100
12) Fluorene	15.510	166	10400	0.353	ng/ul	100
14) Pentachlorophenol	16.875	266	623	0.296	ng/ul	99
15) Phenanthrene	17.251	178	16985	0.348	ng/ul	99
16) Anthracene	17.344	178	13132	0.347	ng/ul	99
19) Fluoranthene	19.276	202	18010	0.366	ng/ul	97
20) Pyrene	19.638	202	18355	0.361	ng/ul	95
21) Benzo(a)anthracene	21.401	228	12452	0.343	ng/ul	99
22) Chrysene	21.451	228	16945	0.354	ng/ul	99
24) Benzo(b)fluoranthene	23.047	252	15450	0.377	ng/ul	98
25) Benzo(k)fluoranthene	23.093	252	15186	0.351	ng/ul	99
26) Benzo(a)pyrene	23.652	252	12774	0.342	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.179	276	14641	0.328	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.196	278	11511	0.326	ng/ul	96
29) Benzo(g,h,i)perylene	26.917	276	13924	0.350	ng/ul	99

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

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