

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062422\
 Data File : BN020425.D
 Acq On : 25 Jun 2022 08:59
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4380

Quant Time: Jun 27 03:42:00 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 24 00:17:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.750	152	3016	0.400	ng/u1	0.00
4) Naphthalene-d8	10.535	136	10115	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.377	164	6932	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.120	188	15868	0.400	ng/u1	0.00
17) Chrysene-d12	21.309	240	14168	0.400	ng/u1	0.00
23) Perylene-d12	23.592	264	11716	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.206	96	1474	0.422	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.124	152	7281	0.439	ng/u1	0.00
18) Fluoranthene-d10	19.151	212	19139	0.410	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.239	88	1425	0.399	ng/u1	94
5) Naphthalene	10.584	128	11980	0.384	ng/u1	99
7) 2-Methylnaphthalene	12.201	142	8156	0.390	ng/u1	99
8) 1-Methylnaphthalene	12.421	142	8770	0.443	ng/u1	100
10) Acenaphthylene	14.095	152	12702	0.387	ng/u1	100
11) Acenaphthene	14.442	153	9901	0.376	ng/u1	99
12) Fluorene	15.427	166	11790	0.378	ng/u1	99
14) Pentachlorophenol	16.787	266	1053	0.329	ng/u1	98
15) Phenanthrene	17.158	178	20415	0.370	ng/u1	100
16) Anthracene	17.251	178	18146	0.381	ng/u1	100
19) Fluoranthene	19.179	202	23377	0.361	ng/u1	99
20) Pyrene	19.541	202	25243	0.376	ng/u1	99
21) Benzo(a)anthracene	21.295	228	19582	0.385	ng/u1	99
22) Chrysene	21.347	228	21334	0.377	ng/u1	100
24) Benzo(b)fluoranthene	22.905	252	19407	0.358	ng/u1	99
25) Benzo(k)fluoranthene	22.946	252	19773	0.361	ng/u1	99
26) Benzo(a)pyrene	23.490	252	18433	0.374	ng/u1	97
27) Indeno(1,2,3-cd)pyrene	25.930	276	18553	0.378	ng/u1#	97
28) Dibenzo(a,h)anthracene	25.943	278	15014	0.379	ng/u1	100
29) Benzo(g,h,i)perylene	26.641	276	15897	0.378	ng/u1	99

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

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