

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN033122\
 Data File : BN019280.D
 Acq On : 01 Apr 2022 09:38
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDCCC0.4

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/04/2022
 Supervised By : Jagrut Upadhyay 04/04/2022

Quant Time: Apr 01 16:40:17 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN033122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 31 12:41:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.833	152	5216	0.400	ng	0.00
7) Naphthalene-d8	10.637	136	14323	0.400	ng	# 0.00
13) Acenaphthene-d10	14.474	164	9045	0.400	ng	0.00
19) Phenanthrene-d10	17.215	188	18373	0.400	ng	# 0.00
29) Chrysene-d12	21.404	240	13554	0.400	ng	# 0.00
35) Perylene-d12	23.761	264	11323	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.392	112	4528	0.397	ng	0.00
5) Phenol-d6	6.995	99	4489	0.371	ng	0.00
8) Nitrobenzene-d5	8.993	82	3551	0.345	ng	0.00
11) 2-Methylnaphthalene-d10	12.231	152	9380	0.412	ng	0.00
14) 2,4,6-Tribromophenol	15.964	330	1412	0.334	ng	0.00
15) 2-Fluorobiphenyl	13.105	172	14110	0.389	ng	0.00
27) Fluoranthene-d10	19.248	212	21571	0.391	ng	0.00
31) Terphenyl-d14	19.847	244	12164	0.403	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.225	88	2460	0.355	ng	90
3) n-Nitrosodimethylamine	3.557	42	2379	0.329	ng	96
6) bis(2-Chloroethyl)ether	7.248	93	5567	0.386	ng	96
9) Naphthalene	10.680	128	16575	0.399	ng	99
10) Hexachlorobutadiene	10.979	225	3889	0.401	ng	# 99
12) 2-Methylnaphthalene	12.303	142	10748	0.401	ng	99
16) Acenaphthylene	14.196	152	15071	0.362	ng	99
17) Acenaphthene	14.538	154	11480	0.387	ng	98
18) Fluorene	15.521	166	14541	0.392	ng	99
20) 4,6-Dinitro-2-methylph...	15.618	198	634	0.292	ng	# 75
21) 4-Bromophenyl-phenylether	16.405	248	4503	0.382	ng	# 80
22) Hexachlorobenzene	16.528	284	6123	0.399	ng	99
23) Atrazine	16.672	200	2764	0.342	ng	98
24) Pentachlorophenol	16.877	266	1557	0.352	ng	98
25) Phenanthrene	17.256	178	22767	0.394	ng	100
26) Anthracene	17.349	178	17403	0.364	ng	100
28) Fluoranthene	19.278	202	26689	0.387	ng	99
30) Pyrene	19.640	202	27133	0.405	ng	99
32) Benzo(a)anthracene	21.386	228	14872	0.350	ng	99
33) Chrysene	21.440	228	21306	0.392	ng	100
34) Bis(2-ethylhexyl)phtha...	21.315	149	9811	0.346	ng	99
36) Indeno(1,2,3-cd)pyrene	26.208	276	15695m	0.339	ng	
37) Benzo(b)fluoranthene	23.045	252	15594	0.371	ng	96
38) Benzo(k)fluoranthene	23.092	252	16735m	0.369	ng	
39) Benzo(a)pyrene	23.659	252	14715	0.385	ng	# 88
40) Dibenzo(a,h)anthracene	26.226	278	10972	0.325	ng	# 92
41) Benzo(g,h,i)perylene	26.945	276	15506	0.329	ng	99

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN033122\
 Data File : BN019280.D
 Acq On : 01 Apr 2022 09:38
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDCCC0.4

Quant Time: Apr 01 16:40:17 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN033122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 31 12:41:03 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 04/04/2022
 Supervised By : Jagrut Upadhyay 04/04/2022

