

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062722\
 Data File : BN020532.D
 Acq On : 28 Jun 2022 14:07
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDCCC0.4EC

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 06/29/2022
 Supervised By : Jagrut Upadhyay 06/30/2022

Quant Time: Jun 28 23:50:15 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN062622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 27 02:03:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.739	152	3097	0.400	ng	0.00
7) Naphthalene-d8	10.520	136	9756	0.400	ng	# 0.00
13) Acenaphthene-d10	14.367	164	6006	0.400	ng	0.00
19) Phenanthrene-d10	17.113	188	12573	0.400	ng	0.00
29) Chrysene-d12	21.306	240	9783	0.400	ng	# 0.00
35) Perylene-d12	23.586	264	7364	0.400	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.341	112	3220	0.375	ng	0.00
5) Phenol-d6	6.944	99	4014	0.391	ng	0.00
8) Nitrobenzene-d5	8.897	82	2841	0.359	ng	0.00
11) 2-Methylnaphthalene-d10	12.117	152	6393	0.382	ng	0.00
14) 2,4,6-Tribromophenol	15.861	330	797	0.352	ng	0.00
15) 2-Fluorobiphenyl	12.988	172	9304	0.375	ng	-0.01
27) Fluoranthene-d10	19.143	212	13677	0.366	ng	0.00
31) Terphenyl-d14	19.746	244	9729	0.416	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.232	88	1455	0.358	ng	98
3) n-Nitrosodimethylamine	3.564	42	1209	0.327	ng	# 92
6) bis(2-Chloroethyl)ether	7.168	93	3296	0.373	ng	99
9) Naphthalene	10.573	128	11217	0.385	ng	99
10) Hexachlorobutadiene	10.861	225	1978	0.373	ng	# 100
12) 2-Methylnaphthalene	12.189	142	7461	0.385	ng	99
16) Acenaphthylene	14.089	152	10673	0.370	ng	99
17) Acenaphthene	14.431	154	7483	0.381	ng	97
18) Fluorene	15.415	166	9661	0.377	ng	100
20) 4,6-Dinitro-2-methylph...	15.522	198	403	0.389	ng	# 64
21) 4-Bromophenyl-phenylether	16.302	248	2703	0.370	ng	95
22) Hexachlorobenzene	16.426	284	3042	0.376	ng	99
23) Atrazine	16.579	200	1896	0.329	ng	98
24) Pentachlorophenol	16.774	266	780	0.484	ng	94
25) Phenanthrene	17.154	178	15769	0.372	ng	100
26) Anthracene	17.246	178	13141	0.360	ng	99
28) Fluoranthene	19.173	202	17149	0.371	ng	100
30) Pyrene	19.535	202	17482	0.422	ng	99
32) Benzo(a)anthracene	21.288	228	12882	0.369	ng	98
33) Chrysene	21.342	228	15096	0.390	ng	99
34) Bis(2-ethylhexyl)phtha...	21.225	149	7039	0.357	ng	99
36) Indeno(1,2,3-cd)pyrene	25.922	276	10739	0.327	ng	98
37) Benzo(b)fluoranthene	22.896	252	11295	0.398	ng	96
38) Benzo(k)fluoranthene	22.943	252	12387m	0.410	ng	
39) Benzo(a)pyrene	23.486	252	9291	0.378	ng	95
40) Dibenzo(a,h)anthracene	25.933	278	8465	0.322	ng	97
41) Benzo(g,h,i)perylene	26.632	276	9508	0.331	ng	100

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

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