

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN030525\
 Data File : BN036524.D
 Acq On : 05 Mar 2025 13:31
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDCCC0.4

Manual Integrations
 APPROVED

Reviewed By :Anahy Claudio 03/06/2025
 Supervised By :Jagrut Upadhyay 03/06/2025

Quant Time: Mar 05 16:09:24 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN021025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 05 16:08:57 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.732	152	2348	0.400	ng	0.00
7) Naphthalene-d8	10.509	136	5891	0.400	ng	0.00
13) Acenaphthene-d10	14.366	164	3887	0.400	ng	0.00
19) Phenanthrene-d10	17.111	188	7928	0.400	ng	0.00
29) Chrysene-d12	21.295	240	7302	0.400	ng	0.00
35) Perylene-d12	23.551	264	7292	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.319	112	2256	0.406	ng	0.00
5) Phenol-d6	6.901	99	2710	0.416	ng	0.00
8) Nitrobenzene-d5	8.875	82	2360	0.406	ng	0.00
11) 2-Methylnaphthalene-d10	12.106	152	3483	0.385	ng	0.00
14) 2,4,6-Tribromophenol	15.858	330	613	0.318	ng	0.00
15) 2-Fluorobiphenyl	12.988	172	8237	0.564	ng	0.00
27) Fluoranthene-d10	19.141	212	8803	0.399	ng	0.00
31) Terphenyl-d14	19.745	244	6133	0.393	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.247	88	1024	0.399	ng	92
3) n-Nitrosodimethylamine	3.550	42	1908	0.428	ng	# 98
6) bis(2-Chloroethyl)ether	7.154	93	2868	0.421	ng	98
9) Naphthalene	10.562	128	6578	0.387	ng	100
10) Hexachlorobutadiene	10.850	225	1524	0.368	ng	# 99
12) 2-Methylnaphthalene	12.182	142	4220	0.379	ng	100
16) Acenaphthylene	14.078	152	6474	0.377	ng	100
17) Acenaphthene	14.430	154	4257	0.371	ng	98
18) Fluorene	15.414	166	5822	0.357	ng	100
20) 4,6-Dinitro-2-methylph...	15.499	198	494	0.318	ng	94
21) 4-Bromophenyl-phenylether	16.305	248	1776	0.376	ng	93
22) Hexachlorobenzene	16.416	284	2169	0.371	ng	98
23) Atrazine	16.578	200	1528	0.387	ng	95
24) Pentachlorophenol	16.764	266	1040	0.375	ng	98
25) Phenanthrene	17.149	178	9019	0.394	ng	99
26) Anthracene	17.235	178	8093	0.400	ng	99
28) Fluoranthene	19.174	202	11196	0.398	ng	98
30) Pyrene	19.536	202	11526	0.410	ng	99
32) Benzo(a)anthracene	21.277	228	9909	0.412	ng	98
33) Chrysene	21.331	228	10684	0.411	ng	98
34) Bis(2-ethylhexyl)phtha...	21.214	149	7086	0.473	ng	100
36) Indeno(1,2,3-cd)pyrene	25.838	276	9728	0.382	ng	100
37) Benzo(b)fluoranthene	22.873	252	10291m	0.429	ng	
38) Benzo(k)fluoranthene	22.917	252	9771	0.395	ng	99
39) Benzo(a)pyrene	23.455	252	8631	0.412	ng	# 91
40) Dibenzo(a,h)anthracene	25.852	278	7487	0.372	ng	99
41) Benzo(g,h,i)perylene	26.528	276	8456	0.371	ng	98

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

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