Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061325\
Data File : BN037259.D
Acq On : 14 Jun 2025 13:23
Operator : RC/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SSTDCCC0.4

Quant Time: Jun 16 02:42:36 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN061325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 13 18:43:34 2025
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	1044	0.400	ng	0.00
7) Naphthalene-d8	10.351	136	2562	0.400	ng	#-0.01
13) Acenaphthene-d10	14.224	164	1268	0.400	ng	0.00
19) Phenanthrene-d10	16.971	188	2318	0.400	ng	0.00
29) Chrysene-d12	21.171	240	1702	0.400	ng	# 0.00
35) Perylene-d12	23.360	264	1789	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.170	112	1026	0.400	ng	0.00
5) Phenol-d6	6.759	99	1096	0.406	ng	0.00
8) Nitrobenzene-d5	8.728	82	1069	0.422	ng	0.00
11) 2-Methylnaphthalene-d10	11.950	152	1407	0.409	ng	0.00
14) 2,4,6-Tribromophenol	15.718	330	230	0.437	ng	-0.01
15) 2-Fluorobiphenyl	12.848	172	2271	0.426	ng	0.00
27) Fluoranthene-d10	19.012	212	2382	0.393	ng	0.00
31) Terphenyl-d14	19.621	244	1544	0.401	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.104	88	562	0.392	ng	92
3) n-Nitrosodimethylamine	3.422	42	1426	0.437	ng	# 97
6) bis(2-Chloroethyl)ether	7.012	93	1026	0.424	ng	94
9) Naphthalene	10.404	128	2969	0.400	ng	100
10) Hexachlorobutadiene	10.693	225	761	0.422	ng	# 97
12) 2-Methylnaphthalene	12.026	142	1778	0.394	ng	97
16) Acenaphthylene	13.946	152	2462	0.396	ng	99
17) Acenaphthene	14.288	154	1581	0.394	ng	98
18) Fluorene	15.282	166	1985	0.385	ng	100
20) 4,6-Dinitro-2-methylph...	15.378	198	199	0.443	ng	90
21) 4-Bromophenyl-phenylether	16.177	248	578	0.383	ng	88
22) Hexachlorobenzene	16.289	284	670	0.383	ng	94
23) Atrazine	16.450	200	489	0.363	ng	# 95
24) Pentachlorophenol	16.636	266	367	0.428	ng	94
25) Phenanthrene	17.009	178	2788	0.379	ng	98
26) Anthracene	17.108	178	2496	0.371	ng	99
28) Fluoranthene	19.045	202	3112	0.362	ng	98
30) Pyrene	19.407	202	3140	0.392	ng	100
32) Benzo(a)anthracene	21.162	228	2322	0.404	ng	99
33) Chrysene	21.206	228	2833	0.396	ng	99
34) Bis(2-ethylhexyl)phtha...	21.099	149	1690	0.395	ng	98
36) Indeno(1,2,3-cd)pyrene	25.547	276	3038	0.421	ng	99
37) Benzo(b)fluoranthene	22.708	252	2598	0.397	ng	99
38) Benzo(k)fluoranthene	22.752	252	2852	0.378	ng	99
39) Benzo(a)pyrene	23.266	252	2372	0.403	ng	96
40) Dibenzo(a,h)anthracene	25.561	278	2198	0.401	ng	97
41) Benzo(g,h,i)perylene	26.216	276	2747	0.411	ng	97

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

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