

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080621\
 Data File : BN015816.D
 Acq On : 06 Aug 2021 20:06
 Operator : CG/JU
 Sample : SSTDICC0.8
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDICC0.8

Quant Time: Aug 07 04:13:41 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN080621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 07 04:10:55 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	8922	0.400	ng	0.00
7) Naphthalene-d8	10.736	136	48347	0.400	ng	0.00
13) Acenaphthene-d10	14.561	164	26107	0.400	ng	0.00
19) Phenanthrene-d10	17.310	188	48553	0.400	ng	0.00
29) Chrysene-d12	21.504	240	44733	0.400	ng	0.00
35) Perylene-d12	23.933	264	45147	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.485	112	20464	0.938	ng	0.00
5) Phenol-d6	7.080	99	28280	1.073	ng	0.00
8) Nitrobenzene-d5	9.076	82	27299	0.879	ng	-0.01
11) 2-Methylnaphthalene-d10	12.318	152	59462	0.819	ng	0.00
14) 2,4,6-Tribromophenol	16.056	330	10333	1.092	ng	0.00
15) 2-Fluorobiphenyl	13.190	172	74550	0.716	ng	0.00
27) Fluoranthene-d10	19.346	212	108749	0.861	ng	0.00
31) Terphenyl-d14	19.942	244	106508	0.962	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.419	88	3037	1.051	ng	97
3) n-Nitrosodimethylamine	3.770	42	6941	1.168	ng	# 98
6) bis(2-Chloroethyl)ether	7.357	93	20511	0.857	ng	99
9) Naphthalene	10.787	128	100375	0.792	ng	99
10) Hexachlorobutadiene	11.078	225	18378	0.831	ng	# 99
12) 2-Methylnaphthalene	12.394	142	67597	0.829	ng	99
16) Acenaphthylene	14.281	152	95856	0.891	ng	100
17) Acenaphthene	14.624	154	60209	0.802	ng	99
18) Fluorene	15.614	166	73117	0.813	ng	99
20) 4,6-Dinitro-2-methylph...	15.677	198	2991	1.016	ng	# 88
21) 4-Bromophenyl-phenylether	16.506	248	20776	0.728	ng	99
22) Hexachlorobenzene	16.616	284	25250	0.770	ng	99
23) Atrazine	16.762	200	20898	1.191	ng	99
24) Pentachlorophenol	16.957	266	9243	0.911	ng	99
25) Phenanthrene	17.346	178	108741	0.753	ng	100
26) Anthracene	17.444	178	106048	0.874	ng	100
28) Fluoranthene	19.371	202	123060	0.819	ng	100
30) Pyrene	19.733	202	125419	0.840	ng	100
32) Benzo(a)anthracene	21.482	228	121031	0.885	ng	100
33) Chrysene	21.535	228	113109	0.765	ng	99
34) Bis(2-ethylhexyl)phtha...	21.419	149	85874	1.567	ng	99
36) Indeno(1,2,3-cd)pyrene	26.456	276	136838	0.856	ng	# 74
37) Benzo(b)fluoranthene	23.190	252	124868	0.796	ng	98
38) Benzo(k)fluoranthene	23.238	252	120566	0.743	ng	99
39) Benzo(a)pyrene	23.823	252	110769	0.850	ng	98
40) Dibenzo(a,h)anthracene	26.480	278	112340	0.839	ng	99
41) Benzo(g,h,i)perylene	27.233	276	117731	0.842	ng	99

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

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