

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101821\
 Data File : BN016987.D
 Acq On : 18 Oct 2021 13:54
 Operator : CG/JU
 Sample : SSTDICC3.2
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDICC3.2

Quant Time: Oct 18 14:24:29 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN101821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 18 13:18:40 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.607	152	19503	0.40	ng	0.00
7) Naphthalene-d8	10.370	136	82936	0.40	ng	0.00
13) Acenaphthene-d10	14.232	164	45807	0.40	ng	0.00
19) Phenanthrene-d10	16.982	188	86391	0.40	ng	0.00
29) Chrysene-d12	21.186	240	86856	0.40	ng	0.00
35) Perylene-d12	23.403	264	82850	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.209	112	138374	2.85	ng	0.00
5) Phenol-d6	6.786	99	163218	3.03	ng	0.00
8) Nitrobenzene-d5	8.747	82	193172	3.66	ng	0.00
11) 2-Methylnaphthalene-d10	11.964	152	376019	3.20	ng	0.00
14) 2,4,6-Tribromophenol	15.729	330	76255	3.70	ng	0.00
15) 2-Fluorobiphenyl	12.849	172	531288	3.05	ng	0.00
27) Fluoranthene-d10	19.019	212	728262	3.32	ng	0.00
31) Terphenyl-d14	19.630	244	611134	3.17	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.199	88	39303	3.12	ng	96
3) n-Nitrosodimethylamine	3.522	42	51421	3.23	ng	# 99
6) bis(2-Chloroethyl)ether	7.044	93	166685	3.08	ng	99
9) Naphthalene	10.420	128	665675	2.85	ng	100
10) Hexachlorobutadiene	10.712	225	132379	3.14	ng	# 100
12) 2-Methylnaphthalene	12.039	142	433076	3.11	ng	100
16) Acenaphthylene	13.953	152	648069	3.42	ng	100
17) Acenaphthene	14.295	154	400748	3.13	ng	98
18) Fluorene	15.285	166	493044	3.22	ng	99
20) 4,6-Dinitro-2-methylph...	15.374	198	49177	3.93	ng	# 80
21) 4-Bromophenyl-phenylether	16.179	248	168597	3.26	ng	99
22) Hexachlorobenzene	16.288	284	182847	3.08	ng	99
23) Atrazine	16.459	200	134538	3.75	ng	98
24) Pentachlorophenol	16.641	266	77345	4.08	ng	100
25) Phenanthrene	17.019	178	758553	3.08	ng	100
26) Anthracene	17.116	178	718872	3.39	ng	100
28) Fluoranthene	19.049	202	855125	3.17	ng	100
30) Pyrene	19.411	202	875833	3.17	ng	100
32) Benzo(a)anthracene	21.165	228	898738	3.39	ng	99
33) Chrysene	21.218	228	862455	3.06	ng	99
34) Bis(2-ethylhexyl)phtha...	21.122	149	494942	3.97	ng	99
36) Indeno(1,2,3-cd)pyrene	25.649	276	945700	3.09	ng	100
37) Benzo(b)fluoranthene	22.739	252	892419	3.25	ng	100
38) Benzo(k)fluoranthene	22.784	252	866858	3.19	ng	99
39) Benzo(a)pyrene	23.308	252	811631	3.35	ng	99
40) Dibenzo(a,h)anthracene	25.669	278	768284	3.10	ng	99
41) Benzo(g,h,i)perylene	26.330	276	828461	3.06	ng	100

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

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