

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102722\
 Data File : BN022342.D
 Acq On : 26 Oct 2022 19:35
 Operator : CG/JU
 Sample : SSTDICC0.8
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDICC0.8

Quant Time: Oct 27 06:52:56 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN102722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 27 06:48:35 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.942	152	3177	0.400	ng	0.00
7) Naphthalene-d8	10.773	136	9718	0.400	ng	0.00
13) Acenaphthene-d10	14.610	164	5481	0.400	ng	0.00
19) Phenanthrene-d10	17.362	188	11555	0.400	ng	0.00
29) Chrysene-d12	21.573	240	9552	0.400	ng	0.00
35) Perylene-d12	24.037	264	8282	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.472	112	6329	0.860	ng	0.00
5) Phenol-d6	7.104	99	7901	0.827	ng	0.00
8) Nitrobenzene-d5	9.119	82	6227	0.797	ng	0.00
11) 2-Methylnaphthalene-d10	12.365	152	11219	0.628	ng	0.00
14) 2,4,6-Tribromophenol	16.109	330	1694	0.810	ng	0.00
15) 2-Fluorobiphenyl	13.231	172	17401	0.727	ng	-0.01
27) Fluoranthene-d10	19.407	212	23743	0.773	ng	0.00
31) Terphenyl-d14	20.002	244	18422	0.959	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.291	88	3415	0.831	ng	99
3) n-Nitrosodimethylamine	3.631	42	3651	0.817	ng	97
6) bis(2-Chloroethyl)ether	7.364	93	7821	0.775	ng	99
9) Naphthalene	10.816	128	21873	0.745	ng	99
10) Hexachlorobutadiene	11.104	225	3748	0.737	ng	# 99
12) 2-Methylnaphthalene	12.073	142	4140	0.952	ng	# 89
16) Acenaphthylene	14.333	152	19318	0.741	ng	99
17) Acenaphthene	14.675	154	13486	0.736	ng	99
18) Fluorene	15.658	166	18755	0.848	ng	100
20) 4,6-Dinitro-2-methylph...	15.749	198	1262	0.766	ng	94
21) 4-Bromophenyl-phenylether	16.556	248	5139	0.730	ng	94
22) Hexachlorobenzene	16.667	284	5946	0.679	ng	100
23) Atrazine	16.829	200	4206	0.805	ng	98
24) Pentachlorophenol	17.027	266	2277	0.848	ng	99
25) Phenanthrene	17.412	178	29109	0.733	ng	100
26) Anthracene	17.499	178	24458	0.767	ng	100
28) Fluoranthene	19.437	202	31981	0.751	ng	100
30) Pyrene	19.803	202	32667	0.779	ng	100
32) Benzo(a)anthracene	21.555	228	27378	0.762	ng	99
33) Chrysene	21.609	228	29443	0.739	ng	100
34) Bis(2-ethylhexyl)phtha...	21.475	149	18607	1.080	ng	98
36) Indeno(1,2,3-cd)pyrene	26.616	276	31675	0.757	ng	100
37) Benzo(b)fluoranthene	23.280	252	28118	0.716	ng	94
38) Benzo(k)fluoranthene	23.330	252	27841	0.712	ng	96
39) Benzo(a)pyrene	23.926	252	23244	0.776	ng	# 91
40) Dibenzo(a,h)anthracene	26.637	278	24743	0.741	ng	98
41) Benzo(g,h,i)perylene	27.411	276	26052	0.723	ng	99

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

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