

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN081019\
 Data File : BN007047.D
 Acq On : 10 Aug 2019 13:06
 Operator : HP/JU
 Sample : SSTDICC0.8
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDICC0.8

Quant Time: Aug 10 21:34:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN081019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 10 21:17:06 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.43	152	6387	0.40	ng	0.00
6) Naphthalene-d8	10.19	136	23207	0.40	ng	0.00
12) Acenaphthene-d10	14.08	164	14131	0.40	ng	0.00
18) Phenanthrene-d10	16.82	188	63780	0.40	ng	0.00
26) Chrysene-d12	21.06	240	73865	0.40	ng	0.00
32) Perylene-d12	23.17	264	74288	0.40	ng	0.00

System Monitoring Compounds						
3) 2-Fluorophenol	5.06	112	11608	0.70	ng	0.00
4) Phenol-d6	6.61	99	14199	0.70	ng	0.00
7) Nitrobenzene-d5	8.56	82	13567	0.75	ng	0.00
10) 2-Methylnaphthalene-d10	11.79	152	31840	0.71	ng	0.00
13) 2,4,6-Tribromophenol	15.57	330	7248	1.09	ng	0.00
14) 2-Fluorobiphenyl	12.71	172	16677	0.71	ng	0.00
24) Fluoranthene-d10	18.87	212	166731	0.71	ng	0.00
28) Terphenyl-d14	19.50	244	145498	0.68	ng	0.00

Target Compounds				Qvalue		
2) n-Nitrosodimethylamine	3.35	42	3108	0.850	ng	# 98
5) bis(2-Chloroethyl)ether	6.87	93	12830	0.712	ng	100
8) Naphthalene	10.24	128	48251	0.707	ng	99
9) Hexachlorobutadiene	10.55	225	10292	0.701	ng	# 100
11) 2-Methylnaphthalene	11.87	142	34015	0.715	ng	97
15) Acenaphthylene	13.79	152	55261	0.744	ng	100
16) Acenaphthene	14.14	154	35515	0.739	ng	98
17) Fluorene	15.14	166	59575	0.855	ng	99
19) 4-Bromophenyl-phenylether	16.04	248	25609	0.603	ng	99
20) Hexachlorobenzene	16.14	284	27029	0.682	ng	100
21) Pentachlorophenol	16.48	266	9888	0.717	ng	99
22) Phenanthrene	16.87	178	138209	0.690	ng	100
23) Anthracene	16.96	178	128276	0.735	ng	100
25) Fluoranthene	18.90	202	182469	0.710	ng	100
27) Pyrene	19.27	202	182224	0.653	ng	100
29) Benzo(a)anthracene	21.03	228	193700	0.693	ng	100
30) Chrysene	21.09	228	185807	0.672	ng	100
31) Indeno(1,2,3-cd)pyrene	25.25	276	203408	0.688	ng	100
33) Benzo(b)fluoranthene	22.55	252	185486	0.717	ng	98
34) Benzo(k)fluoranthene	22.59	252	180198	0.697	ng	98
35) Benzo(a)pyrene	23.08	252	164269	0.704	ng	97
36) Dibenzo(a,h)anthracene	25.27	278	166033	0.714	ng	99
37) Benzo(g,h,i)perylene	25.88	276	168730	0.695	ng	98

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

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