

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN083019\
 Data File : BN007520.D
 Acq On : 30 Aug 2019 13:03
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDCCC0.4

Quant Time: Aug 31 01:58:32 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN082319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 29 12:18:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	6859	0.40	ng	-0.02
7) Naphthalene-d8	10.14	136	31010	0.40	ng	-0.03
13) Acenaphthene-d10	14.03	164	19635	0.40	ng	-0.02
19) Phenanthrene-d10	16.79	188	47308	0.40	ng	0.00
27) Chrysene-d12	21.02	240	41883	0.40	ng	0.00
34) Perylene-d12	23.12	264	38962	0.40	ng	-0.01

System Monitoring Compounds						
4) 2-Fluorophenol	5.02	112	8639	0.36	ng	-0.02
5) Phenol-d6	6.58	99	10885	0.36	ng	0.00
8) Nitrobenzene-d5	8.52	82	10271	0.38	ng	-0.01
11) 2-Methylnaphthalene-d10	11.75	152	21168	0.41	ng	-0.01
14) 2,4,6-Tribromophenol	15.54	330	4710	0.46	ng	-0.02
15) 2-Fluorobiphenyl	12.65	172	30599	0.39	ng	0.00
25) Fluoranthene-d10	18.83	212	58870	0.39	ng	-0.01
29) Terphenyl-d14	19.46	244	47449	0.43	ng	-0.01

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.00	88	4995	0.438	ng	# 71
3) n-Nitrosodimethylamine	3.32	42	1735	0.328	ng	# 67
6) bis(2-Chloroethyl)ether	6.83	93	10416	0.420	ng	95
9) Naphthalene	10.19	128	34791	0.393	ng	# 90
10) Hexachlorobutadiene	10.49	225	6512	0.359	ng	# 99
12) 2-Methylnaphthalene	11.82	142	25295	0.419	ng	95
16) Acenaphthylene	13.75	152	43184	0.372	ng	99
17) Acenaphthene	14.10	154	26406	0.388	ng	98
18) Fluorene	15.09	166	34840	0.405	ng	100
20) 4-Bromophenyl-phenylether	16.01	248	2100	0.113	ng	# 81
21) Hexachlorobenzene	16.10	284	12381	0.390	ng	95
22) Pentachlorophenol	16.45	266	5953	0.604	ng	# 63
23) Phenanthrene	16.83	178	58392	0.410	ng	99
24) Anthracene	16.92	178	56478	0.395	ng	99
26) Fluoranthene	18.86	202	71869	0.393	ng	99
28) Pyrene	19.23	202	73811	0.438	ng	99
30) Benzo(a)anthracene	20.99	228	66080	0.402	ng	99
31) Chrysene	21.05	228	62738	0.395	ng	99
32) Bis(2-ethylhexyl)phthalate	20.96	149	56020	0.507	ng	99
33) Indeno(1,2,3-cd)pyrene	25.18	276	60135	0.314	ng	96
35) Benzo(b)fluoranthene	22.50	252	59052	0.418	ng	97
36) Benzo(k)fluoranthene	22.54	252	60119	0.431	ng	# 69
37) Benzo(a)pyrene	23.03	252	54406	0.405	ng	98
38) Dibenzo(a,h)anthracene	25.19	278	47489	0.354	ng	93
39) Benzo(g,h,i)perylene	25.80	276	48442	0.363	ng	99

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

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