

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN032819\
 Data File : BN005019.D
 Acq On : 28 Mar 2019 19:13
 Operator : JU/SJ
 Sample : SSTDICC0.8
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDICC0.8

Quant Time: Mar 29 03:23:55 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN032819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 29 03:18:03 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	2727	0.40	ng	0.00
7) Naphthalene-d8	10.44	136	11749	0.40	ng	0.00
13) Acenaphthene-d10	14.30	164	7228	0.40	ng	0.00
19) Phenanthrene-d10	17.06	188	17326	0.40	ng	0.00
27) Chrysene-d12	21.26	240	21650	0.40	ng	0.00
34) Perylene-d12	23.50	264	20985	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.25	112	5590	0.75	ng	0.00
5) Phenol-d6	6.84	99	6589	0.71	ng	0.00
8) Nitrobenzene-d5	8.82	82	5505	0.71	ng	-0.01
11) 2-Methylnaphthalene-d10	12.04	152	15092	0.74	ng	0.00
14) 2,4,6-Tribromophenol	15.81	330	3162	0.63	ng	0.00
15) 2-Fluorobiphenyl	12.92	172	22483	0.75	ng	0.00
25) Fluoranthene-d10	19.10	212	40388	0.70	ng	0.00
29) Terphenyl-d14	19.69	244	34623	0.78	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.18	88	2294	0.740	ng	97
3) n-Nitrosodimethylamine	3.52	42	1409	0.769	ng	# 93
6) bis(2-Chloroethyl)ether	7.10	93	6382	0.766	ng	99
9) Naphthalene	10.49	128	24390	0.740	ng	99
10) Hexachlorobutadiene	10.76	225	4475	0.712	ng	# 99
12) 2-Methylnaphthalene	12.11	142	17698	0.723	ng	99
16) Acenaphthylene	14.02	152	27781	0.746	ng	100
17) Acenaphthene	14.37	154	17767	0.729	ng	100
18) Fluorene	15.36	166	24052	0.707	ng	100
20) 4-Bromophenyl-phenylether	16.25	248	8174	0.722	ng	96
21) Hexachlorobenzene	16.36	284	9019	0.696	ng	100
22) Pentachlorophenol	16.73	266	1678	0.662	ng	94
23) Phenanthrene	17.10	178	39612	0.708	ng	100
24) Anthracene	17.20	178	35423	0.709	ng	100
26) Fluoranthene	19.13	202	48600	0.676	ng	100
28) Pyrene	19.49	202	49208	0.745	ng	100
30) Benzo(a)anthracene	21.25	228	48072	0.677	ng	99
31) Chrysene	21.30	228	52053	0.696	ng	99
32) Bis(2-ethylhexyl)phthalate	21.15	149	18473	0.651	ng	99
33) Indeno(1,2,3-cd)pyrene	25.74	276	60652	0.561	ng	100
35) Benzo(b)fluoranthene	22.82	252	55476	0.731	ng	98
36) Benzo(k)fluoranthene	22.86	252	53100	0.711	ng	98
37) Benzo(a)pyrene	23.40	252	49138	0.706	ng	97
38) Dibenzo(a,h)anthracene	25.75	278	49595	0.615	ng	98
39) Benzo(g,h,i)perylene	26.43	276	49693	0.605	ng	99

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

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