

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061019\
 Data File : BE099885.D
 Acq On : 10 Jun 2019 10:18
 Operator : JU/SJ
 Sample : SSTDIC0.1
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDIC0.1

Manual Integrations
 APPROVED

mohammad
 6/14/2019 8:38:56 AM

Quant Time: Jun 10 14:45:27 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061019.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Mon Jun 10 14:31:35 2019

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	821	0.40	ng	0.00
7) Naphthalene-d8	10.61	136	2978	0.40	ng	0.01
13) Acenaphthene-d10	14.47	164	2153	0.40	ng	0.00
19) Phenanthrene-d10	17.21	188	6495	0.40	ng	0.00
27) Chrysene-d12	21.40	240	9777	0.40	ng	0.00
34) Perylene-d12	23.93	264	10541	0.40	ng	0.00

System Monitoring Compounds

4) 2-Fluorophenol	5.36	112	205	0.07	ng	0.00
5) Phenol-d6	6.98	99	252	0.07	ng	0.00
8) Nitrobenzene-d5	8.99	82	238	0.08	ng	0.03
11) 2-Methylnaphthalene-d10	12.22	152	467	0.09	ng	0.01
14) 2,4,6-Tribromophenol	15.98	330	121	0.06	ng	0.01
15) 2-Fluorobiphenyl	13.12	172	777	0.10	ng	0.01
25) Fluoranthene-d10	19.26	212	8202	0.08	ng	0.00
29) Terphenyl-d14	19.85	244	1624	0.09	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.28	88	132	0.108	ng	# 92
6) bis(2-Chloroethyl)ether	7.23	93	192	0.077	ng	# 95
9) Naphthalene	10.65	128	3071	0.095	ng	# 92
10) Hexachlorobutadiene	10.93	225	237	0.100	ng	# 91
12) 2-Methylnaphthalene	12.30	142	467	0.085	ng	# 89
16) Acenaphthylene	14.20	152	984	0.092	ng	# 93
17) Acenaphthene	14.54	154	529	0.091	ng	# 99
18) Fluorene	15.51	166	799	0.089	ng	# 95
20) 4-Bromophenyl-phenylether	16.41	248	326	0.088	ng	# 80
21) Hexachlorobenzene	16.52	284	371	0.102	ng	# 94
23) Phenanthrene	17.27	178	1434	0.089	ng	# 99
24) Anthracene	17.36	178	1136	0.074	ng	# 98
26) Fluoranthene	19.28	202	2312	0.081	ng	# 99
28) Pyrene	19.65	202	2457	0.095	ng	# 99
30) Benzo(a)anthracene	21.39	228	2641m	0.085	ng	# 97
31) Chrysene	21.44	228	2841	0.093	ng	# 97
32) Bis(2-ethylhexyl)phthalate	21.29	149	4978	0.098	ng	# 100
33) Indeno(1,2,3-cd)pyrene	26.62	276	3391	0.091	ng	# 99
35) Benzo(b)fluoranthene	23.15	252	2678	0.091	ng	# 78
36) Benzo(k)fluoranthene	23.20	252	2742	0.094	ng	# 82
37) Benzo(a)pyrene	23.82	252	2778	0.098	ng	# 65
38) Dibenzo(a,h)anthracene	26.65	278	2611	0.090	ng	# 81
39) Benzo(g,h,i)perylene	27.44	276	2767	0.094	ng	# 86

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

