

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE060719\
 Data File : BE099846.D
 Acq On : 7 Jun 2019 11:05
 Operator : JU/SJ
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICC0.1

Quant Time: Jun 07 14:50:58 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE060719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 07 14:29:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	558	0.40	ng	0.01
7) Naphthalene-d8	10.61	136	2167	0.40	ng	0.01
13) Acenaphthene-d10	14.48	164	1913	0.40	ng	0.01
19) Phenanthrene-d10	17.23	188	6690	0.40	ng	0.01
27) Chrysene-d12	21.41	240	10358	0.40	ng	0.01
34) Perylene-d12	23.94	264	11309	0.40	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.36	112	234	0.17	ng	0.00
5) Phenol-d6	6.98	99	247	0.14	ng	0.01
8) Nitrobenzene-d5	8.98	82	172	0.09	ng	0.01
11) 2-Methylnaphthalene-d10	12.22	152	377	0.10	ng	0.01
14) 2,4,6-Tribromophenol	15.98	330	158	0.14	ng	0.01
15) 2-Fluorobiphenyl	13.12	172	592	0.08	ng	0.01
25) Fluoranthene-d10	19.26	212	9489	0.10	ng	0.00
29) Terphenyl-d14	19.86	244	2005	0.12	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.27	88	134	0.166	ng	# 76
6) bis(2-Chloroethyl)ether	7.23	93	187	0.122	ng	# 85
9) Naphthalene	10.67	128	2315	0.100	ng	# 91
10) Hexachlorobutadiene	10.94	225	161	0.091	ng	# 98
12) 2-Methylnaphthalene	12.30	142	393	0.107	ng	# 88
16) Acenaphthylene	14.20	152	825	0.089	ng	# 94
17) Acenaphthene	14.54	154	476	0.094	ng	# 96
18) Fluorene	15.53	166	722	0.097	ng	# 95
20) 4-Bromophenyl-phenylether	16.41	248	369	0.099	ng	# 79
21) Hexachlorobenzene	16.53	284	345	0.084	ng	# 98
23) Phenanthrene	17.27	178	1537	0.096	ng	# 98
24) Anthracene	17.36	178	1429	0.099	ng	# 97
26) Fluoranthene	19.29	202	2755	0.098	ng	# 100
28) Pyrene	19.66	202	2848	0.114	ng	# 97
30) Benzo(a)anthracene	21.39	228	3239	0.113	ng	# 83
31) Chrysene	21.45	228	3033	0.096	ng	# 86
32) Bis(2-ethylhexyl)phthalate	21.31	149	6557	0.200	ng	# 98
33) Indeno(1,2,3-cd)pyrene	26.64	276	3618	0.096	ng	# 100
35) Benzo(b)fluoranthene	23.16	252	3147	0.101	ng	# 56
36) Benzo(k)fluoranthene	23.22	252	2971	0.089	ng	# 59
37) Benzo(a)pyrene	23.83	252	2993	0.099	ng	# 43
38) Dibenzo(a,h)anthracene	26.66	278	2853	0.092	ng	# 49
39) Benzo(g,h,i)perylene	27.48	276	2998	0.092	ng	# 74

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

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