

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE052219\
 Data File : BE099713.D
 Acq On : 22 May 2019 10:06
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC0.4

Manual Integrations
 APPROVED

mohammad
 5/24/2019 8:47:48 AM

Quant Time: May 22 11:54:20 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE052019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 20 19:29:40 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	2095	0.40	ng	0.00
7) Naphthalene-d8	10.61	136	8237	0.40	ng	0.00
13) Acenaphthene-d10	14.49	164	5909	0.40	ng	0.00
19) Phenanthrene-d10	17.23	188	17458	0.40	ng	0.00
27) Chrysene-d12	21.41	240	23551	0.40	ng	0.00
34) Perylene-d12	23.96	264	27834	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.37	112	1992	0.38	ng	0.00
5) Phenol-d6	6.97	99	2715	0.39	ng	0.00
8) Nitrobenzene-d5	8.98	82	2540	0.36	ng	0.00
11) 2-Methylnaphthalene-d10	12.22	152	5100	0.36	ng	0.00
14) 2,4,6-Tribromophenol	15.98	330	1243	0.41	ng	0.00
15) 2-Fluorobiphenyl	13.10	172	7916	0.35	ng	0.00
25) Fluoranthene-d10	19.26	212	79109	0.33	ng	0.00
29) Terphenyl-d14	19.86	244	15980	0.38	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.30	88	1010	0.326	ng	93
3) n-Nitrosodimethylamine	3.63	42	555	0.324	ng	# 96
6) bis(2-Chloroethyl)ether	7.24	93	2448	0.393	ng	93
9) Naphthalene	10.67	128	31902	0.362	ng	100
10) Hexachlorobutadiene	10.94	225	2268	0.350	ng	99
12) 2-Methylnaphthalene	12.30	142	5220	0.363	ng	99
16) Acenaphthylene	14.20	152	9814	0.356	ng	100
17) Acenaphthene	14.54	154	5601	0.362	ng	97
18) Fluorene	15.53	166	8425	0.370	ng	98
20) 4-Bromophenyl-phenylether	16.43	248	3622	0.360	ng	94
21) Hexachlorobenzene	16.53	284	3888	0.377	ng	99
22) Pentachlorophenol	16.91	266	754	0.418	ng	95
23) Phenanthrene	17.27	178	15734	0.369	ng	100
24) Anthracene	17.36	178	13903	0.359	ng	100
26) Fluoranthene	19.29	202	23093	0.342	ng	99
28) Pyrene	19.65	202	23692	0.400	ng	100
30) Benzo(a)anthracene	21.40	228	27111	0.405	ng	99
31) Chrysene	21.45	228	28168	0.394	ng	99
32) Bis(2-ethylhexyl)phthalate	21.32	149	34544	0.448	ng	99
33) Indeno(1,2,3-cd)pyrene	26.64	276	34412	0.398	ng	99
35) Benzo(b)fluoranthene	23.17	252	28385m	0.372	ng	
36) Benzo(k)fluoranthene	23.22	252	28484	0.358	ng	100
37) Benzo(a)pyrene	23.84	252	26833	0.366	ng	# 98
38) Dibenzo(a,h)anthracene	26.67	278	28174	0.366	ng	# 98
39) Benzo(g,h,i)perylene	27.48	276	27814	0.357	ng	99

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

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