

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE050118\
 Data File : BE096267.D
 Acq On : 1 May 2018 21:06
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC0.4EC

Manual Integrations
 APPROVED

Sohil
 5/2/2018 4:36:27 AM

Quant Time: May 02 03:47:05 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE041118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 30 14:27:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	801	0.40	ng	0.00
7) Naphthalene-d8	11.19	136	3097	0.40	ng	0.00
13) Acenaphthene-d10	14.94	164	1772	0.40	ng	0.00
19) Phenanthrene-d10	17.65	188	4223	0.40	ng	0.00
27) Chrysene-d12	21.74	240	4076	0.40	ng	0.00
34) Perylene-d12	24.35	264	3969	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.88	112	986	0.40	ng	0.01
5) Phenol-d6	7.52	99	1390	0.41	ng	0.01
8) Nitrobenzene-d5	9.54	82	1472m	0.41	ng	0.00
11) 2-Methylnaphthalene-d10	12.73	152	1904	0.39	ng	0.00
14) 2,4,6-Tribromophenol	16.41	330	350	0.41	ng	0.00
15) 2-Fluorobiphenyl	13.57	172	2830	0.37	ng	0.00
25) Fluoranthene-d10	19.64	212	20287	0.35	ng	0.00
29) Terphenyl-d14	20.19	244	3644	0.40	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.61	88	404	0.327	ng	87
3) n-Nitrosodimethylamine	3.96	42	629	0.415	ng	# 78
6) bis(2-Chloroethyl)ether	7.76	93	1186	0.394	ng	90
9) Naphthalene	11.24	128	12307	0.377	ng	99
10) Hexachlorobutadiene	11.50	225	458	0.316	ng	95
12) 2-Methylnaphthalene	12.81	142	2082	0.385	ng	96
16) Acenaphthylene	14.67	152	3611	0.382	ng	99
17) Acenaphthene	15.00	154	2160	0.380	ng	95
18) Fluorene	15.97	166	3354	0.477	ng	# 72
20) 4-Bromophenyl-phenylether	16.83	248	957	0.381	ng	# 71
21) Hexachlorobenzene	16.96	284	934	0.374	ng	# 81
22) Pentachlorophenol	17.31	266	407	0.489	ng	# 80
23) Phenanthrene	17.70	178	4359	0.367	ng	99
24) Anthracene	17.78	178	4316	0.381	ng	# 94
26) Fluoranthene	19.67	202	5492	0.352	ng	97
28) Pyrene	20.00	202	341	0.042	ng	# 79
30) Benzo(a)anthracene	21.73	228	4986	0.381	ng	98
31) Chrysene	21.79	228	4856	0.385	ng	95
32) Bis(2-ethylhexyl)phthalate	21.60	149	15526	0.460	ng	100
33) Indeno(1,2,3-cd)pyrene	27.13	276	4978	0.407	ng	94
35) Benzo(b)fluoranthene	23.54	252	4475	0.376	ng	94
36) Benzo(k)fluoranthene	23.59	252	4579	0.374	ng	# 83
37) Benzo(a)pyrene	24.23	252	4331	0.378	ng	# 84
38) Dibenzo(a,h)anthracene	27.14	278	4163	0.413	ng	# 90
39) Benzo(g,h,i)perylene	28.00	276	4179	0.392	ng	# 91

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

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