

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE092118\
 Data File : BE097642.D
 Acq On : 22 Sep 2018 10:35
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC0.4EC

Manual Integrations
 APPROVED

Sohil
 9/24/2018 8:31:08 AM

Quant Time: Sep 24 02:59:51 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE091918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 20 10:54:43 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.37	152	870	0.40	ng	-0.01
7) Naphthalene-d8	10.13	136	4378	0.40	ng	0.00
13) Acenaphthene-d10	14.00	164	3083	0.40	ng	0.00
19) Phenanthrene-d10	16.75	188	9530	0.40	ng	0.00
27) Chrysene-d12	20.97	240	10232	0.40	ng	0.00
34) Perylene-d12	23.20	264	9960	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.05	112	904	0.35	ng	0.00
5) Phenol-d6	6.61	99	1306m	0.39	ng	0.00
8) Nitrobenzene-d5	8.53	82	1133	0.34	ng	0.00
11) 2-Methylnaphthalene-d10	11.72	152	2560	0.42	ng	0.00
14) 2,4,6-Tribromophenol	15.52	330	515	0.44	ng	0.00
15) 2-Fluorobiphenyl	12.62	172	3932	0.36	ng	0.00
25) Fluoranthene-d10	18.80	212	38906	0.32	ng	0.00
29) Terphenyl-d14	19.42	244	7730	0.41	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.07	88	580	0.389	ng	Qvalue # 83
3) n-Nitrosodimethylamine	3.38	42	438m	0.349	ng	
6) bis(2-Chloroethyl)ether	6.83	93	1104	0.357	ng	76
9) Naphthalene	10.17	128	15825	0.401	ng	99
10) Hexachlorobutadiene	10.46	225	704	0.387	ng	97
12) 2-Methylnaphthalene	11.80	142	2570	0.418	ng	98
16) Acenaphthylene	13.71	152	4883	0.401	ng	99
17) Acenaphthene	14.06	154	3266	0.400	ng	96
18) Fluorene	15.05	166	4468	0.419	ng	# 79
20) 4-Bromophenyl-phenylether	15.96	248	1688	0.393	ng	# 87
21) Hexachlorobenzene	16.07	284	1857	0.393	ng	# 85
22) Pentachlorophenol	16.44	266	447	0.341	ng	# 77
23) Phenanthrene	16.80	178	8820	0.390	ng	99
24) Anthracene	16.89	178	7281	0.364	ng	96
26) Fluoranthene	18.83	202	10304	0.328	ng	98
28) Pyrene	19.20	202	10459	0.449	ng	100
30) Benzo(a)anthracene	20.95	228	9204	0.357	ng	97
31) Chrysene	21.00	228	11145	0.403	ng	99
32) Bis(2-ethylhexyl)phthalate	20.91	149	11884	0.285	ng	99
33) Indeno(1,2,3-cd)pyrene	25.50	276	12853	0.386	ng	# 93
35) Benzo(b)fluoranthene	22.53	252	9416	0.369	ng	# 90
36) Benzo(k)fluoranthene	22.58	252	11714	0.402	ng	95
37) Benzo(a)pyrene	23.11	252	9456	0.383	ng	92
38) Dibenzo(a,h)anthracene	25.52	278	10514	0.393	ng	96
39) Benzo(g,h,i)perylene	26.20	276	10202	0.392	ng	# 88

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

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