

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE090216\
 Data File : BE092348.D
 Acq On : 2 Sep 2016 18:03
 Operator : UM/SJ
 Sample : SSTDICC005
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICC005

Manual Integrations
 APPROVED

sohil
 9/6/2016 1:15:20 PM

Quant Time: Sep 03 00:17:39 2016
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 02 18:37:51 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.16	152	8867	5.00	ng	-0.02
7) Naphthalene-d8	10.95	136	44418	5.00	ng	0.00
12) Acenaphthene-d10	14.82	164	27581	5.00	ng	0.00
18) Phenanthrene-d10	17.58	188	62412	5.00	ng	0.02
24) Chrysene-d12	21.75	240	96900	5.00	ng	0.00
31) Perylene-d12	24.11	264	81941	5.00	ng	0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.68	112	11887	5.88	ng	-0.01
5) Phenol-d6	7.33	99	17608	6.53	ng	-0.02
8) Nitrobenzene-d5	9.34	82	17352	5.66	ng	-0.02
13) 2,4,6-Tribromophenol	0.00	330	0d	0.00	ng	
14) 2-Fluorobiphenyl	13.43	172	43048	5.07	ng	-0.01
26) Terphenyl-d14	20.17	244	66036	4.48	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.41	88	5718	4.38	ng	Qvalue 93
9) Naphthalene	10.99	128	50211	5.02	ng	# 95
10) Hexachlorobutadiene	11.30	225	7700	4.89	ng	99
11) 2-Methylnaphthalene	12.62	142	34244	5.31	ng	95
15) Acenaphthylene	14.53	152	248033	5.29	ng	100
16) Acenaphthene	14.88	154	37652	4.83	ng	96
17) Fluorene	15.86	166	49847	5.19	ng	99
19) Hexachlorobenzene	16.89	284	15392	5.16	ng	# 64
21) Phenanthrene	17.60	178	84338	4.73	ng	# 94
22) Anthracene	17.69	178	89528	5.97	ng	98
23) Fluoranthene	19.61	202	400537	5.29	ng	99
25) Pyrene	19.97	202	444539	4.62	ng	100
27) Benzo(a)anthracene	21.72	228	531156m	5.00	ng	
28) Chrysene	21.77	228	457470m	4.06	ng	
30) Indeno(1,2,3-cd)pyrene	26.60	276	503340	4.43	ng	99
32) Benzo(b)fluoranthene	23.37	252	108314	5.16	ng	94
33) Benzo(k)fluoranthene	23.43	252	118744m	5.40	ng	
34) Benzo(a)pyrene	24.00	252	109807	5.36	ng	# 94
35) Dibenzo(a,h)anthracene	26.62	278	103998	5.53	ng	95
36) Benzo(g,h,i)perylene	27.37	276	422110	5.26	ng	97

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

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