

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE021517\
 Data File : BE092735.D
 Acq On : 15 Feb 2017 16:13
 Operator : SJ/MA
 Sample : SSTDICC010
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

mohammad
 2/16/2017 11:47:20 AM

Quant Time: Feb 15 17:03:30 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE021517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 15 16:44:55 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.10	152	10095	5.00	ng	-0.02
7) Naphthalene-d8	10.88	136	50166	5.00	ng	-0.02
12) Acenaphthene-d10	14.76	164	31384	5.00	ng	0.00
18) Phenanthrene-d10	17.51	188	67325	5.00	ng	0.00
24) Chrysene-d12	21.70	240	71105	5.00	ng	0.00
31) Perylene-d12	24.03	264	70631	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	0.00	112	0d	0.00	ng	
5) Phenol-d6	0.00	99	0d	0.00	ng	
8) Nitrobenzene-d5	0.00	82	0d	0.00	ng	
13) 2,4,6-Tribromophenol	0.00	330	0d	0.00	ng	
14) 2-Fluorobiphenyl	13.37	172	79185	10.40	ng	-0.02
26) Terphenyl-d14	20.12	244	105071	11.85	ng	-0.02
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.34	88	12524	8.58	ng	92
3) n-Nitrosodimethylamine	3.69	42	19424	13.59	ng	98
9) Naphthalene	10.93	128	106745	9.81	ng	94
10) Hexachlorobutadiene	11.22	225	15841	9.91	ng	99
11) 2-Methylnaphthalene	12.55	142	73770	10.94	ng	98
15) Acenaphthylene	14.47	152	509365	10.87	ng	100
16) Acenaphthene	14.81	154	72857	10.68	ng	93
17) Fluorene	15.81	166	94086	10.75	ng	100
19) Hexachlorobenzene	16.84	284	22697	8.58	ng	# 39
21) Phenanthrene	17.56	178	147904m	9.15	ng	
22) Anthracene	17.65	178	148492	10.09	ng	98
23) Fluoranthene	19.54	202	690784	10.08	ng	98
25) Pyrene	19.92	202	768350	11.57	ng	99
27) Benzo(a)anthracene	21.68	228	716769	10.45	ng	# 76
28) Chrysene	21.72	228	842080m	10.34	ng	
30) Indeno(1,2,3-cd)pyrene	26.48	276	859494	9.23	ng	98
32) Benzo(b)fluoranthene	23.30	252	187815	9.33	ng	94
33) Benzo(k)fluoranthene	23.35	252	198065	9.30	ng	93
34) Benzo(a)pyrene	23.91	252	186320	10.30	ng	94
35) Dibenzo(a,h)anthracene	26.50	278	179630	9.30	ng	96
36) Benzo(g,h,i)perylene	27.23	276	745874	9.12	ng	98

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

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