

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE090115\
 Data File : BE090673.D
 Acq On : 1 Sep 2015 19:19
 Operator : UM/IZ
 Sample : SSTDICCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICCC001

Manual Integrations
 APPROVED

MMDadoda
 9/2/2015 4:07:41 PM

Quant Time: Sep 02 01:42:00 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 02 01:28:15 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	48112	5.00	ng	0.00
6) Naphthalene-d8	10.31	136	204950	5.00	ng	0.00
12) Acenaphthene-d10	14.17	164	99119	5.00	ng	0.00
18) Phenanthrene-d10	16.91	188	249198	5.00	ng	0.00
25) Chrysene-d12	21.07	240	313353	5.00	ng	0.00
32) Perylene-d12	23.36	264	274127	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.19	112	9391	0.70	ng	0.00
5) Phenol-d6	6.75	99	12724	0.65	ng	0.00
7) Nitrobenzene-d5	8.71	82	13515	0.84	ng	0.00
13) 2,4,6-Tribromophenol	15.68	330	2393	0.65	ng	0.00
14) 2-Fluorobiphenyl	12.80	172	27032	0.92	ng	0.00
27) Terphenyl-d14	19.54	244	33113	0.61	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.16	88	7619	1.18	ng	100
3) n-Nitrosodimethylamine	3.45	42	9411	1.11	ng	100
8) Nitrobenzene	8.75	77	14390	0.86	ng	100
9) Naphthalene	10.36	128	44501	0.96	ng	100
10) Hexachlorobutadiene	10.65	225	7233	1.05	ng	100
11) 2-Methylnaphthalene	11.99	142	23432	0.79	ng	100
15) Acenaphthylene	13.89	152	152097	0.86	ng	100
16) Acenaphthene	14.23	154	26766	0.98	ng	100
17) Fluorene	15.22	166	32713	0.98	ng	100
19) 4-Bromophenyl-phenylether	16.13	248	8167	0.82	ng	100
20) Hexachlorobenzene	16.23	284	45615	1.01	ng	99
21) Pentachlorophenol	16.60	266	2356m	0.62	ng	
22) Phenanthrene	16.96	178	61488	0.95	ng	100
23) Anthracene	17.05	178	49570	0.84	ng	100
24) Fluoranthene	18.97	202	60252	0.88	ng	100
26) Pyrene	19.33	202	61450	0.71	ng	100
28) Benzo(a)anthracene	21.06	228	68536	0.84	ng	100
29) Chrysene	21.11	228	79246	0.95	ng	100
30) Bis(2-ethylhexyl)phthalate	21.01	149	15888	0.37	ng	100
31) Indeno(1,2,3-cd)pyrene	25.73	276	68317	0.87	ng	100
33) Benzo(b)fluoranthene	22.67	252	57889	0.90	ng	100
34) Benzo(k)fluoranthene	22.71	252	68224	0.94	ng	100
35) Benzo(a)pyrene	23.26	252	49681	0.83	ng	100
36) Dibenzo(a,h)anthracene	25.75	278	53749	0.92	ng	100
37) Benzo(g,h,i)perylene	26.46	276	57908	0.91	ng	100

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

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