

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102414\
 Data File : BE088212.D
 Acq On : 24 Oct 2014 14:54
 Operator : TP/JJ
 Sample : SSTDICC002
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICC002

Manual Integrations
 APPROVED

mohammad
 10/28/2014 12:22:18 PM

Quant Time: Oct 27 12:51:56 2014
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE102414.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 27 12:38:23 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.11	152	56013	5.00	ng	0.00
3) Naphthalene-d8	8.68	136	216263	5.00	ng	0.00
7) Acenaphthene-d10	10.83	164	98675	5.00	ng	0.00
12) Phenanthrene-d10	12.94	188	175743	5.00	ng	0.00
16) Chrysene-d12	18.35	240	98346	5.00	ng	0.00
22) Perylene-d12	21.40	264	87841	5.00	ng	0.00
System Monitoring Compounds						
4) Nitrobenzene-d5	7.80	82	29926	2.25	ng	0.00
8) 2-Fluorobiphenyl	10.01	172	56931	2.07	ng	0.00
18) Terphenyl-d14	16.11	244	34954	2.17	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.40	88	16131	2.05	ng	100
5) Naphthalene	8.70	128	90554	2.06	ng	99
6) 2-Methylnaphthalene	9.55	142	56417	2.13	ng	99
9) Acenaphthylene	10.64	152	317980	2.16	ng	100
10) Acenaphthene	10.86	154	46655	2.00	ng	98
11) Fluorene	11.52	166	50854	2.14	ng	100
13) Phenanthrene	12.97	178	334724	2.31	ng	97
14) Anthracene	13.06	178	264596	2.11	ng	100
15) Fluoranthene	15.21	202	55055	2.14	ng	100
17) Pyrene	15.67	202	59336	2.20	ng	100
19) Benzo(a)anthracene	18.33	228	161265	2.08	ng	100
20) Chrysene	18.40	228	178587	2.03	ng	100
21) Indeno(1,2,3-cd)pyrene	23.58	276	154689m	1.99	ng	
23) Benzo(b)fluoranthene	20.64	252	34063	2.37	ng	99
24) Benzo(k)fluoranthene	20.70	252	47473	2.09	ng	99
25) Benzo(a)pyrene	21.28	252	33359	2.35	ng	99
26) Dibenzo(a,h)anthracene	23.68	278	30326	2.26	ng	97
27) Benzo(g,h,i)perylene	24.20	276	152272	2.16	ng	99

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

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