

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102114\
 Data File : BE088160.D
 Acq On : 22 Oct 2014 1:54
 Operator : TP/JJ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Manual Integrations
 APPROVED

mohammad
 10/22/2014 1:55:26 PM

Quant Time: Oct 22 04:25:20 2014
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE102014.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 21 02:11:14 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.14	152	31672	5.00	ng	0.00
3) Naphthalene-d8	8.70	136	122757	5.00	ng	0.00
7) Acenaphthene-d10	10.85	164	57106	5.00	ng	0.00
12) Phenanthrene-d10	12.96	188	81744	5.00	ng	0.00
16) Chrysene-d12	18.38	240	54054	5.00	ng	0.00
22) Perylene-d12	21.44	264	46314	5.00	ng	0.00
System Monitoring Compounds						
4) Nitrobenzene-d5	7.83	82	8905	1.01	ng	0.00
8) 2-Fluorobiphenyl	10.04	172	16480	1.04	ng	0.00
18) Terphenyl-d14	16.14	244	10316	1.19	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.34	88	4651	1.02	ng	99
5) Naphthalene	8.72	128	26273	1.04	ng	100
6) 2-Methylnaphthalene	9.58	142	16012	1.06	ng	99
9) Acenaphthylene	10.67	152	90835	1.04	ng	100
10) Acenaphthene	10.88	154	13904	1.04	ng	99
11) Fluorene	11.55	166	14543	1.02	ng	99
13) Phenanthrene	13.00	178	81402m	1.18	ng	
14) Anthracene	13.09	178	76111	1.19	ng	99
15) Fluoranthene	15.25	202	16223	1.04	ng	100
17) Pyrene	15.70	202	17618	1.20	ng	100
19) Benzo(a)anthracene	18.36	228	44288	0.92	ng	99
20) Chrysene	18.44	228	49016	1.04	ng	98
21) Indeno(1,2,3-cd)pyrene	23.64	276	44598m	0.85	ng	
23) Benzo(b)fluoranthene	20.68	252	9634	0.87	ng	99
24) Benzo(k)fluoranthene	20.74	252	11673	0.99	ng	99
25) Benzo(a)pyrene	21.32	252	8967	0.88	ng	100
26) Dibenzo(a,h)anthracene	23.73	278	8895	0.89	ng	100
27) Benzo(g,h,i)perylene	24.26	276	43367	0.99	ng	99

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

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