

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE103114\
 Data File : BE088454.D
 Acq On : 1 Nov 2014 17:02
 Operator : TP/JJ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Manual Integrations
 APPROVED

mohammad
 11/4/2014 9:07:21 AM

Quant Time: Nov 02 04:33:37 2014
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE103114.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 01 00:42:07 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.08	152	27190	5.00	ng	0.00
3) Naphthalene-d8	8.64	136	103453	5.00	ng	-0.01
7) Acenaphthene-d10	10.79	164	46786	5.00	ng	0.00
12) Phenanthrene-d10	12.89	188	89397	5.00	ng	0.00
16) Chrysene-d12	18.30	240	62524	5.00	ng	-0.01
22) Perylene-d12	21.36	264	58787	5.00	ng	-0.01
System Monitoring Compounds						
4) Nitrobenzene-d5	7.77	82	7831	1.03	ng	0.00
8) 2-Fluorobiphenyl	9.98	172	13549	1.03	ng	0.00
18) Terphenyl-d14	16.06	244	10852	0.94	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	2.31	88	3771	1.06	ng	99
5) Naphthalene	8.66	128	22061	1.03	ng	100
6) 2-Methylnaphthalene	9.52	142	13424	1.01	ng	94
9) Acenaphthylene	10.61	152	77367	1.02	ng	99
10) Acenaphthene	10.83	154	11517	1.02	ng	99
11) Fluorene	11.48	166	12560	1.05	ng	100
13) Phenanthrene	12.93	178	96871	1.19	ng	99
14) Anthracene	13.02	178	72800	1.11	ng	99
15) Fluoranthene	15.17	202	16995	1.29	ng	100
17) Pyrene	15.62	202	18280	0.90	ng	100
19) Benzo(a)anthracene	18.27	228	61839	1.05	ng	97
20) Chrysene	18.35	228	59457	1.03	ng	100
21) Indeno(1,2,3-cd)pyrene	23.52	276	61265m	1.14	ng	
23) Benzo(b)fluoranthene	20.60	252	13885	1.08	ng	100
24) Benzo(k)fluoranthene	20.66	252	14816	1.04	ng	100
25) Benzo(a)pyrene	21.24	252	13358	1.08	ng	98
26) Dibenzo(a,h)anthracene	23.62	278	12163	1.06	ng	99
27) Benzo(g,h,i)perylene	24.14	276	55158	1.05	ng	99

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

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