

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE103114\
 Data File : BE088427.D
 Acq On : 31 Oct 2014 17:33
 Operator : TP/JJ
 Sample : SSTDICC002
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICC002

Manual Integrations
 APPROVED

mohammad
 11/4/2014 9:05:46 AM

Quant Time: Oct 31 23:40:33 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 00:20:37 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.09	152	39904	5.00	ng	0.00
3) Naphthalene-d8	8.65	136	153929	5.00	ng	0.00
7) Acenaphthene-d10	10.79	164	72762	5.00	ng	0.00
12) Phenanthrene-d10	12.90	188	138058	5.00	ng	0.00
16) Chrysene-d12	18.31	240	66047	5.00	ng	0.00
22) Perylene-d12	21.37	264	57510	5.00	ng	0.00
System Monitoring Compounds						
4) Nitrobenzene-d5	7.77	82	23633	2.10	ng	0.00
8) 2-Fluorobiphenyl	9.98	172	41990	2.03	ng	0.00
18) Terphenyl-d14	16.07	244	25196	2.07	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.37	88	10796	2.07	ng	99
5) Naphthalene	8.68	128	65233	2.02	ng	97
6) 2-Methylnaphthalene	9.53	142	41118	2.06	ng	98
9) Acenaphthylene	10.62	152	240580	2.01	ng	99
10) Acenaphthene	10.84	154	34839	1.96	ng	99
11) Fluorene	11.49	166	38417	2.04	ng	99
13) Phenanthrene	12.94	178	274456	1.96	ng	99
14) Anthracene	13.02	178	205313	2.01	ng	99
15) Fluoranthene	15.18	202	42535	2.09	ng	100
17) Pyrene	15.63	202	44882	2.07	ng	100
19) Benzo(a)anthracene	18.29	228	125309	2.01	ng	100
20) Chrysene	18.36	228	124733	2.04	ng	99
21) Indeno(1,2,3-cd)pyrene	23.55	276	116988m	2.01	ng	
23) Benzo(b)fluoranthene	20.61	252	25979	2.07	ng	99
24) Benzo(k)fluoranthene	20.67	252	29223	2.10	ng	100
25) Benzo(a)pyrene	21.26	252	24985	2.08	ng	98
26) Dibenzo(a,h)anthracene	23.64	278	22922	2.05	ng	99
27) Benzo(g,h,i)perylene	24.17	276	103389	2.00	ng	98

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

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