

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE103114\
 Data File : BE088424.D
 Acq On : 31 Oct 2014 15:52
 Operator : TP/JJ
 Sample : SSTDICC0.5
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICC0.5

Manual Integrations
 APPROVED

mohammad
 11/4/2014 9:05:38 AM

Quant Time: Oct 31 23:37:46 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	39470	5.00	ng	0.00
3) Naphthalene-d8	8.65	136	152039	5.00	ng	0.00
7) Acenaphthene-d10	10.79	164	71375	5.00	ng	0.00
12) Phenanthrene-d10	12.90	188	139555	5.00	ng	0.00
16) Chrysene-d12	18.31	240	65396	5.00	ng	0.00
22) Perylene-d12	21.37	264	55964	5.00	ng	0.00
System Monitoring Compounds						
4) Nitrobenzene-d5	7.77	82	5659	0.51	ng	0.00
8) 2-Fluorobiphenyl	9.99	172	10560	0.52	ng	0.00
18) Terphenyl-d14	16.08	244	6151	0.51	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.38	88	2603	0.50	ng	100
5) Naphthalene	8.67	128	16375	0.51	ng	98
6) 2-Methylnaphthalene	9.53	142	10181	0.52	ng	98
9) Acenaphthylene	10.62	152	59491	0.51	ng	99
10) Acenaphthene	10.84	154	9107	0.52	ng	98
11) Fluorene	11.49	166	9842	0.53	ng	100
13) Phenanthrene	12.94	178	71240	0.50	ng	97
14) Anthracene	13.02	178	52630	0.51	ng	99
15) Fluoranthene	15.18	202	10461	0.51	ng	100
17) Pyrene	15.64	202	11062	0.52	ng	100
19) Benzo(a)anthracene	18.29	228	30518	0.49	ng	97
20) Chrysene	18.37	228	30801	0.51	ng	97
21) Indeno(1,2,3-cd)pyrene	23.55	276	26798m	0.47	ng	
23) Benzo(b)fluoranthene	20.61	252	6107	0.50	ng	99
24) Benzo(k)fluoranthene	20.67	252	6795	0.50	ng	98
25) Benzo(a)pyrene	21.26	252	5848	0.50	ng	99
26) Dibenzo(a,h)anthracene	23.64	278	5405	0.50	ng	96
27) Benzo(g,h,i)perylene	24.17	276	25107	0.50	ng	98

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

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