

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE032415\
 Data File : BE089838.D
 Acq On : 25 Mar 2015 8:58
 Operator : TP/JJ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001EC

Manual Integrations
 APPROVED

mohammad
 3/25/2015 5:42:50 PM

Quant Time: Mar 25 17:25:36 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIMPAH-BE032415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 25 16:20:31 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	179860	5.00	ng	0.00
4) Naphthalene-d8	10.26	136	707028	5.00	ng	0.00
8) Acenaphthene-d10	14.40	164	322863	5.00	ng	0.01
13) Phenanthrene-d10	17.92	188	552302	5.00	ng	0.00
19) Chrysene-d12	24.26	240	509010	5.00	ng	0.00
25) Perylene-d12	27.41	264	459690	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	8.68	82	6119	0.37	ng	0.00
9) 2-Fluorobiphenyl	12.85	172	10140	0.13	ng	0.00
21) Terphenyl-d14	21.87	244	8462	0.11	ng	0.00
Target Compounds						Qvalue
3) n-Nitrosodimethylamine	3.03	42	3144m	0.09	ng	
6) Naphthalene	10.31	128	15627	0.11	ng	97
7) 2-Methylnaphthalene	11.93	142	8661	0.11	ng	96
10) Acenaphthylene	14.04	152	50175	0.11	ng	97
11) Acenaphthene	14.46	154	7869	0.11	ng	93
12) Fluorene	15.71	166	9513	0.11	ng	92
14) Hexachlorobenzene	16.98	284	2051	0.11	ng	99
16) Phenanthrene	17.97	178	14013	0.12	ng	98
17) Anthracene	18.10	178	11838	0.11	ng	97
18) Fluoranthene	20.82	202	12937	0.12	ng	99
20) Pyrene	21.35	202	13171	0.10	ng	99
22) Benzo(a)anthracene	24.25	228	11961	0.13	ng	98
23) Chrysene	24.32	228	11680	0.11	ng	94
24) Indeno(1,2,3-cd)pyrene	29.38	276	7814m	0.49	ng	
26) Benzo(b)fluoranthene	26.64	252	7013	0.38	ng	95
27) Benzo(k)fluoranthene	26.70	252	10512m	0.11	ng	
28) Benzo(a)pyrene	27.30	252	7439	0.37	ng	97
29) Dibenzo(a,h)anthracene	29.45	278	3926	0.32	ng	96
30) Benzo(g,h,i)perylene	29.84	276	8493	0.34	ng	96

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

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