

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE120914\
 Data File : BE089195.D
 Acq On : 9 Dec 2014 19:50
 Operator : TP/JJ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Manual Integrations
 APPROVED

mohammad
 12/10/2014 4:31:53 PM

Quant Time: Dec 10 01:01:05 2014
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE120514.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 08 00:50:53 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	22171	5.00	ng	-0.02
3) Naphthalene-d8	8.51	136	86401	5.00	ng	-0.02
7) Acenaphthene-d10	10.65	164	40764	5.00	ng	-0.02
12) Phenanthrene-d10	12.72	188	61465	5.00	ng	-0.02
17) Chrysene-d12	18.07	240	41368	5.00	ng	-0.02
23) Perylene-d12	21.11	264	36730	5.00	ng	-0.03
System Monitoring Compounds						
4) Nitrobenzene-d5	7.64	82	6928	1.09	ng	-0.02
8) 2-Fluorobiphenyl	9.85	172	12267	1.09	ng	-0.01
19) Terphenyl-d14	15.84	244	8056	1.09	ng	-0.03
Target Compounds						Qvalue
2) 1,4-Dioxane	2.19	88	2861	1.03	ng	99
5) Naphthalene	8.54	128	18882	1.04	ng	98
6) 2-Methylnaphthalene	9.39	142	12120	1.09	ng	98
9) Acenaphthylene	10.47	152	70884	1.10	ng	100
10) Acenaphthene	10.69	154	10602	1.07	ng	98
11) Fluorene	11.34	166	11312	1.07	ng	100
13) Pentachlorophenol	12.43	266	958	1.00	ng	98
14) Phenanthrene	12.75	178	61763	1.06	ng	100
15) Anthracene	12.84	178	61977	1.07	ng	100
16) Fluoranthene	14.95	202	13090	1.04	ng	100
18) Pyrene	15.40	202	13746	1.08	ng	100
20) Benzo(a)anthracene	18.05	228	40029	1.04	ng	99
21) Chrysene	18.12	228	40519	1.05	ng	98
22) Indeno(1,2,3-cd)pyrene	23.20	276	37025m	1.09	ng	
24) Benzo(b)fluoranthene	20.36	252	8348	1.08	ng	100
25) Benzo(k)fluoranthene	20.42	252	9532	1.04	ng	99
26) Benzo(a)pyrene	21.00	252	8001	1.06	ng	99
27) Dibenzo(a,h)anthracene	23.28	278	7212	1.10	ng	99
28) Benzo(g,h,i)perylene	23.78	276	33207	1.06	ng	99

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

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