

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102114\
 Data File : BE088141.D
 Acq On : 21 Oct 2014 14:10
 Operator : TP/JJ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Manual Integrations
 APPROVED

mohammad
 10/22/2014 1:53:30 PM

Quant Time: Oct 22 04:36:16 2014
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE102014.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 21 02:11:14 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.14	152	38169	5.00	ng	0.00
3) Naphthalene-d8	8.70	136	142442	5.00	ng	0.00
7) Acenaphthene-d10	10.85	164	63619	5.00	ng	0.00
12) Phenanthrene-d10	12.96	188	115323	5.00	ng	0.00
16) Chrysene-d12	18.38	240	84537	5.00	ng	0.00
22) Perylene-d12	21.44	264	75993	5.00	ng	0.00
System Monitoring Compounds						
4) Nitrobenzene-d5	7.82	82	10283	1.00	ng	0.00
8) 2-Fluorobiphenyl	10.03	172	18432	1.05	ng	0.00
18) Terphenyl-d14	16.14	244	13592	1.00	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.39	88	5707	1.04	ng	98
5) Naphthalene	8.72	128	30519	1.04	ng	100
6) 2-Methylnaphthalene	9.58	142	18285	1.04	ng	98
9) Acenaphthylene	10.66	152	101555	1.05	ng	100
10) Acenaphthene	10.88	154	15516	1.04	ng	98
11) Fluorene	11.55	166	16453	1.04	ng	100
13) Phenanthrene	13.00	178	94551m	0.93	ng	
14) Anthracene	13.09	178	89904	1.00	ng	100
15) Fluoranthene	15.25	202	21344	0.97	ng	99
17) Pyrene	15.70	202	22894	1.00	ng	100
19) Benzo(a)anthracene	18.36	228	70622	0.94	ng	98
20) Chrysene	18.43	228	75702	1.03	ng	98
21) Indeno(1,2,3-cd)pyrene	23.63	276	71824m	0.88	ng	
23) Benzo(b)fluoranthene	20.68	252	16627	0.92	ng	99
24) Benzo(k)fluoranthene	20.74	252	19565	1.01	ng	100
25) Benzo(a)pyrene	21.32	252	15328	0.92	ng	99
26) Dibenzo(a,h)anthracene	23.73	278	14604	0.89	ng	100
27) Benzo(g,h,i)perylene	24.26	276	67751	0.94	ng	99

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

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