

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102414\
 Data File : BE088213.D
 Acq On : 24 Oct 2014 15:27
 Operator : TP/JJ
 Sample : SSTDICC005
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICC005

Manual Integrations
 APPROVED

mohammad
 10/28/2014 12:22:21 PM

Quant Time: Oct 27 12:54:05 2014
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE102414.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 27 12:38:23 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.11	152	61708	5.00	ng	0.00
3) Naphthalene-d8	8.67	136	237895	5.00	ng	0.00
7) Acenaphthene-d10	10.82	164	108788	5.00	ng	0.00
12) Phenanthrene-d10	12.94	188	193936	5.00	ng	0.00
16) Chrysene-d12	18.35	240	113340	5.00	ng	0.00
22) Perylene-d12	21.40	264	103914	5.00	ng	0.00
System Monitoring Compounds						
4) Nitrobenzene-d5	7.80	82	87205	5.97	ng	0.00
8) 2-Fluorobiphenyl	10.02	172	148033	4.88	ng	0.00
18) Terphenyl-d14	16.11	244	101790	5.48	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.39	88	43720	5.05	ng	99
5) Naphthalene	8.70	128	236781	4.89	ng	98
6) 2-Methylnaphthalene	9.55	142	147810	5.07	ng	96
9) Acenaphthylene	10.65	152	822782	5.06	ng	100
10) Acenaphthene	10.86	154	118558	4.61	ng	98
11) Fluorene	11.52	166	129887	4.95	ng	100
13) Phenanthrene	12.97	178	884043	5.52	ng	98
14) Anthracene	13.06	178	717609	5.20	ng	100
15) Fluoranthene	15.21	202	157051	5.54	ng	100
17) Pyrene	15.67	202	166775	5.36	ng	100
19) Benzo(a)anthracene	18.32	228	491347	5.51	ng	99
20) Chrysene	18.40	228	498522	4.92	ng	99
21) Indeno(1,2,3-cd)pyrene	23.58	276	494767m	5.53	ng	
23) Benzo(b)fluoranthene	20.64	252	107283	6.32	ng	99
24) Benzo(k)fluoranthene	20.70	252	136662	5.09	ng	99
25) Benzo(a)pyrene	21.28	252	107076	6.38	ng	98
26) Dibenzo(a,h)anthracene	23.68	278	99738	6.28	ng	97
27) Benzo(g,h,i)perylene	24.21	276	469526	5.63	ng	98

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

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