

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102414\
 Data File : BE088210.D
 Acq On : 24 Oct 2014 13:46
 Operator : TP/JJ
 Sample : SSTDICC0.5
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICC0.5

Manual Integrations
 APPROVED

mohammad
 10/28/2014 12:22:11 PM

Quant Time: Oct 27 12:48:41 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	58770	5.00	ng	0.00
3) Naphthalene-d8	8.68	136	224962	5.00	ng	0.00
7) Acenaphthene-d10	10.82	164	102142	5.00	ng	0.00
12) Phenanthrene-d10	12.94	188	178486	5.00	ng	0.00
16) Chrysene-d12	18.34	240	101054	5.00	ng	0.00
22) Perylene-d12	21.40	264	91713	5.00	ng	0.00
System Monitoring Compounds						
4) Nitrobenzene-d5	7.80	82	6513	0.47	ng	0.00
8) 2-Fluorobiphenyl	10.01	172	14994	0.53	ng	0.00
18) Terphenyl-d14	16.11	244	8644	0.52	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.39	88	4177	0.51	ng	100
5) Naphthalene	8.70	128	23938	0.52	ng	99
6) 2-Methylnaphthalene	9.55	142	14462	0.52	ng	100
9) Acenaphthylene	10.64	152	80056	0.52	ng	100
10) Acenaphthene	10.86	154	12689	0.53	ng	99
11) Fluorene	11.52	166	12958	0.53	ng	100
13) Phenanthrene	12.97	178	89924	0.61	ng	94
14) Anthracene	13.06	178	66573	0.52	ng	99
15) Fluoranthene	15.21	202	13711	0.53	ng	99
17) Pyrene	15.67	202	14678	0.53	ng	100
19) Benzo(a)anthracene	18.32	228	39876	0.50	ng	99
20) Chrysene	18.40	228	46692	0.52	ng	98
21) Indeno(1,2,3-cd)pyrene	23.57	276	34683m	0.44	ng	
23) Benzo(b)fluoranthene	20.64	252	7211	0.48	ng	99
24) Benzo(k)fluoranthene	20.70	252	12419	0.52	ng	99
25) Benzo(a)pyrene	21.28	252	7073	0.48	ng	98
26) Dibenzo(a,h)anthracene	23.68	278	6626	0.47	ng	96
27) Benzo(g,h,i)perylene	24.20	276	36956	0.50	ng	98

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

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