

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE103114\
 Data File : BE088452.D
 Acq On : 1 Nov 2014 8:40
 Operator : TP/JJ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001EC

Manual Integrations
 APPROVED

mohammad
 11/4/2014 9:07:32 AM

Quant Time: Nov 02 04:31:42 2014
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE103114.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 01 00:42:07 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.09	152	21559	5.00	ng	0.00
3) Naphthalene-d8	8.65	136	90965	5.00	ng	0.00
7) Acenaphthene-d10	10.80	164	43497	5.00	ng	0.01
12) Phenanthrene-d10	12.91	188	80631	5.00	ng	0.00
16) Chrysene-d12	18.33	240	35892	5.00	ng	0.02
22) Perylene-d12	21.39	264	34779	5.00	ng	0.02
System Monitoring Compounds						
4) Nitrobenzene-d5	7.78	82	42386	6.35	ng	0.00
8) 2-Fluorobiphenyl	9.99	172	122260	9.97	ng	0.00
18) Terphenyl-d14	16.09	244	73955	11.22	ng	0.02
Target Compounds						Qvalue
2) 1,4-Dioxane	2.25	88	2679	0.95	ng	98
5) Naphthalene	8.68	128	18158	0.96	ng	98
6) 2-Methylnaphthalene	9.53	142	11170	0.96	ng	97
9) Acenaphthylene	10.63	152	65558	0.93	ng	99
10) Acenaphthene	10.84	154	9484	0.91	ng	98
11) Fluorene	11.49	166	10081	0.91	ng	99
13) Phenanthrene	12.95	178	69696	0.95	ng	96
14) Anthracene	13.04	178	54298	0.92	ng	100
15) Fluoranthene	15.19	202	12327	1.04	ng	100
17) Pyrene	15.65	202	12551	1.07	ng	99
19) Benzo(a)anthracene	18.31	228	36643	1.08	ng	98
20) Chrysene	18.38	228	36462	1.10	ng	98
21) Indeno(1,2,3-cd)pyrene	23.57	276	42675m	1.38	ng	
23) Benzo(b)fluoranthene	20.63	252	10883	1.43	ng	98
24) Benzo(k)fluoranthene	20.69	252	8623	1.03	ng	97
25) Benzo(a)pyrene	21.27	252	8987	1.23	ng	98
26) Dibenzo(a,h)anthracene	23.66	278	6923	1.02	ng	95
27) Benzo(g,h,i)perylene	24.20	276	38890	1.25	ng	99

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

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