

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102714\
 Data File : BE088327.D
 Acq On : 27 Oct 2014 23:02
 Operator : TP/JJ
 Sample : PB79870BS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB79870BS

Manual Integrations
 APPROVED

mohammad
 10/29/2014 11:11:38 AM

Quant Time: Oct 28 10:24:42 2014
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE102714.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 28 10:16:06 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.11	152	31885	5.00	ng	0.00
3) Naphthalene-d8	8.67	136	122936	5.00	ng	0.00
7) Acenaphthene-d10	10.81	164	57295	5.00	ng	0.00
12) Phenanthrene-d10	12.93	188	109836	5.00	ng	0.00
16) Chrysene-d12	18.35	240	59241	5.00	ng	0.00
22) Perylene-d12	21.40	264	55031	5.00	ng	0.00
System Monitoring Compounds						
4) Nitrobenzene-d5	7.80	82	82540	8.91	ng	0.00
8) 2-Fluorobiphenyl	10.01	172	139848	9.27	ng	0.00
18) Terphenyl-d14	16.11	244	103976	9.39	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.38	88	42966	9.51	ng	99
5) Naphthalene	8.69	128	223705	8.70	ng	100
6) 2-Methylnaphthalene	9.55	142	144113	8.64	ng	97
9) Acenaphthylene	10.64	152	781129	9.46	ng	100
10) Acenaphthene	10.86	154	117657	9.68	ng	99
11) Fluorene	11.52	166	130454	10.10	ng	98
13) Phenanthrene	12.96	178	918759	10.00	ng	98
14) Anthracene	13.06	178	748403	9.92	ng	99
15) Fluoranthene	15.20	202	168517	9.97	ng	99
17) Pyrene	15.67	202	174560	9.78	ng	99
19) Benzo(a)anthracene	18.32	228	550688	9.86	ng	99
20) Chrysene	18.40	228	529451	9.98	ng	98
21) Indeno(1,2,3-cd)pyrene	23.58	276	555481m	9.18	ng	
23) Benzo(b)fluoranthene	20.64	252	124504	9.97	ng	99
24) Benzo(k)fluoranthene	20.70	252	132889	10.25	ng	99
25) Benzo(a)pyrene	21.29	252	122963	10.23	ng	99
26) Dibenzo(a,h)anthracene	23.68	278	112480	9.60	ng	96
27) Benzo(g,h,i)perylene	24.22	276	476544	9.40	ng	98

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

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