

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102214\
 Data File : BE088182.D
 Acq On : 22 Oct 2014 15:27
 Operator : TP/JJ
 Sample : PB79727BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB79727BL

Manual Integrations
 APPROVED

mohammad
 10/28/2014 12:05:42 PM

Quant Time: Oct 27 04:21:25 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	28319	5.00	ng	0.00
3) Naphthalene-d8	8.70	136	112845	5.00	ng	0.00
7) Acenaphthene-d10	10.85	164	58985	5.00	ng	0.00
12) Phenanthrene-d10	12.97	188	88184	5.00	ng	0.00
16) Chrysene-d12	18.39	240	59895	5.00	ng	0.00
22) Perylene-d12	21.45	264	55660	5.00	ng	0.00
System Monitoring Compounds						
4) Nitrobenzene-d5	7.83	82	61167	7.52	ng	0.00
8) 2-Fluorobiphenyl	10.04	172	109794	6.74	ng	0.00
18) Terphenyl-d14	16.15	244	80222	8.32	ng	0.00
Target Compounds						Qvalue
5) Naphthalene	8.72	128	8655	0.37	ng	100
6) 2-Methylnaphthalene	9.58	142	10288	0.74	ng	96
9) Acenaphthylene	10.67	152	82622	0.92	ng	100
10) Acenaphthene	10.88	154	13370	0.96	ng	99
11) Fluorene	11.55	166	15478	1.05	ng	99
13) Phenanthrene	13.01	178	107349	1.49	ng	96
14) Anthracene	13.09	178	86429	1.26	ng	100
15) Fluoranthene	15.25	202	18095	1.08	ng	99
17) Pyrene	15.70	202	19026	1.17	ng	100
19) Benzo(a)anthracene	18.37	228	52326	0.98	ng	99
20) Chrysene	18.44	228	57693	1.10	ng	99
21) Indeno(1,2,3-cd)pyrene	23.64	276	54449m	0.94	ng	
23) Benzo(b)fluoranthene	20.69	252	11462	0.87	ng	99
24) Benzo(k)fluoranthene	20.74	252	13789	0.97	ng	100
25) Benzo(a)pyrene	21.33	252	11258	0.92	ng	99
26) Dibenzo(a,h)anthracene	23.74	278	10703	0.89	ng	100
27) Benzo(g,h,i)perylene	24.27	276	50649	0.96	ng	98

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

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