

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE032415\
 Data File : BE089821.D
 Acq On : 24 Mar 2015 19:32
 Operator : TP/JJ
 Sample : PB82216BSD
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB82216BSD

Manual Integrations
 APPROVED

mohammad
 3/25/2015 5:43:16 PM

Quant Time: Mar 25 03:36:06 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIMPAH-BE032415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 24 18:03:12 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	173302	5.00	ng	0.00
4) Naphthalene-d8	10.26	136	720331	5.00	ng	0.00
8) Acenaphthene-d10	14.40	164	318580	5.00	ng	0.00
13) Phenanthrene-d10	17.92	188	527260	5.00	ng	0.00
19) Chrysene-d12	24.26	240	493412	5.00	ng	0.00
25) Perylene-d12	27.41	264	412392	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	8.68	82	482376	9.16	ng	0.00
9) 2-Fluorobiphenyl	12.85	172	709211	9.51	ng	0.00
21) Terphenyl-d14	21.86	244	658844	9.05	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.55	88	180716	8.12	ng	91
3) n-Nitrosodimethylamine	2.95	42	263764	8.00	ng	# 77
6) Naphthalene	10.31	128	1055583	7.41	ng	96
7) 2-Methylnaphthalene	11.93	142	637842	8.06	ng	99
10) Acenaphthylene	14.04	152	3708577	8.14	ng	99
11) Acenaphthene	14.47	154	534586	7.66	ng	84
12) Fluorene	15.71	166	674606	7.97	ng	93
14) Hexachlorobenzene	16.99	284	139525	7.96	ng	94
15) Pentachlorophenol	17.50	266	111648	13.37	ng	99
16) Phenanthrene	17.97	178	906212	7.82	ng	100
17) Anthracene	18.10	178	920750	8.89	ng	99
18) Fluoranthene	20.82	202	937936	8.98	ng	99
20) Pyrene	21.34	202	987770	7.89	ng	97
22) Benzo(a)anthracene	24.25	228	819735	8.88	ng	97
23) Chrysene	24.32	228	870499	8.64	ng	99
24) Indeno(1,2,3-cd)pyrene	29.39	276	682647m	6.77	ng	
26) Benzo(b)fluoranthene	26.64	252	682785	7.69	ng	98
27) Benzo(k)fluoranthene	26.70	252	904856	10.71	ng	# 95
28) Benzo(a)pyrene	27.30	252	678575	8.14	ng	# 90
29) Dibenzo(a,h)anthracene	29.44	278	574908	7.33	ng	# 78
30) Benzo(g,h,i)perylene	29.83	276	694434	7.98	ng	# 89

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

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