

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE013015\
 Data File : BE089444.D
 Acq On : 30 Jan 2015 18:43
 Operator : TP/IZ
 Sample : PB81495BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB81495BL

Quant Time: Feb 02 00:07:18 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIMPAH-BE013015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 31 05:05:18 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.59	152	73815	5.00	ng	0.00
2) Naphthalene-d8	12.50	136	306656	5.00	ng	0.00
6) Acenaphthene-d10	16.69	164	165033	5.00	ng	-0.01
11) Phenanthrene-d10	19.99	188	276824	5.00	ng	0.00
16) Chrysene-d12	23.84	240	226573	5.00	ng	0.00
22) Perylene-d12	25.56	264	203436	5.00	ng	0.00

System Monitoring Compounds						
3) Nitrobenzene-d5	10.85	82	127114	6.26	ng	0.00
7) 2-Fluorobiphenyl	15.14	172	204661	4.71	ng	0.00
18) Terphenyl-d14	22.48	244	180727	5.33	ng	0.00

Target Compounds				Qvalue		
4) Naphthalene	12.54	128	4381	0.06	ng	97
5) 2-Methylnaphthalene	14.21	142	5797	0.13	ng	99
8) Acenaphthylene	16.34	152	48532	0.16	ng	99
9) Acenaphthene	16.76	154	7935	0.18	ng	95
10) Fluorene	18.02	166	8702	0.18	ng	99
12) Pentachlorophenol	19.66	266	2274	1.66	ng	92
13) Phenanthrene	20.03	178	56837	0.21	ng	98
14) Anthracene	20.12	178	45407	0.18	ng	100
15) Fluoranthene	21.91	202	10730	0.18	ng	99
17) Pyrene	22.23	202	12023	0.20	ng	99
19) Benzo(a)anthracene	23.83	228	32772	0.17	ng	98
20) Chrysene	23.87	228	47479	0.22	ng	97
21) Indeno(1,2,3-cd)pyrene	27.06	276	16865	0.09	ng	99
23) Benzo(b)fluoranthene	25.10	252	3231	0.27	ng	# 88
24) Benzo(k)fluoranthene	25.13	252	6897	0.13	ng	# 94
25) Benzo(a)pyrene	25.48	252	3806	0.22	ng	# 88
26) Dibenzo(a,h)anthracene	27.08	278	2803	0.20	ng	# 83
27) Benzo(g,h,i)perylene	27.51	276	22906	0.14	ng	# 92

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

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