

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE013015\
 Data File : BE089448.D
 Acq On : 30 Jan 2015 21:18
 Operator : TP/IZ
 Sample : G1150-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 FB

Quant Time: Feb 02 00:18:53 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	81087	5.00	ng	0.00
2) Naphthalene-d8	12.50	136	349429	5.00	ng	0.00
6) Acenaphthene-d10	16.69	164	177850	5.00	ng	-0.01
11) Phenanthrene-d10	19.99	188	290914	5.00	ng	0.00
16) Chrysene-d12	23.84	240	246649	5.00	ng	0.00
22) Perylene-d12	25.55	264	213948	5.00	ng	0.00
System Monitoring Compounds						
3) Nitrobenzene-d5	10.85	82	157711	6.82	ng	0.00
7) 2-Fluorobiphenyl	15.13	172	270404	5.77	ng	0.00
18) Terphenyl-d14	22.48	244	260238	7.05	ng	0.00
Target Compounds						Qvalue
4) Naphthalene	12.55	128	10099	0.12	ng	98
5) 2-Methylnaphthalene	14.20	142	6586	0.13	ng	94
8) Acenaphthylene	16.34	152	41977	0.13	ng	99
9) Acenaphthene	16.76	154	6694	0.14	ng	97
10) Fluorene	18.02	166	7493	0.14	ng	99
12) Pentachlorophenol	19.66	266	2625	1.81	ng	90
13) Phenanthrene	20.03	178	47739	0.17	ng	99
14) Anthracene	20.12	178	41060	0.16	ng	99
15) Fluoranthene	21.91	202	10147	0.16	ng	99
17) Pyrene	22.22	202	10666	0.16	ng	99
19) Benzo(a)anthracene	23.83	228	31668	0.15	ng	99
20) Chrysene	23.87	228	42497	0.18	ng	98
21) Indeno(1,2,3-cd)pyrene	27.06	276	18280	0.09	ng	# 88
23) Benzo(b)fluoranthene	25.10	252	3321	0.27	ng	# 91
24) Benzo(k)fluoranthene	25.13	252	8576	0.16	ng	# 93
25) Benzo(a)pyrene	25.48	252	4085	0.22	ng	# 86
26) Dibenzo(a,h)anthracene	27.08	278	2443	0.19	ng	# 87
27) Benzo(g,h,i)perylene	27.51	276	21023	0.12	ng	90

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

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