

Data Path : Z:\HPCHEM1\BNA E\DATA\BE021815\
 Data File : BE089638.D
 Acq On : 19 Feb 2015 15:15
 Operator : TP/IZ
 Sample : MDL-07-S
 Misc : MDL-SOIL-0.1/0.2
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 MDL-07-S

Quant Time: Feb 20 01:17:30 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SIMPAH-BE021815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 18 13:01:39 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.34	152	35006	5.00	ng	0.00
4) Naphthalene-d8	12.24	136	138578	5.00	ng	0.00
8) Acenaphthene-d10	16.42	164	69078	5.00	ng	0.00
13) Phenanthrene-d10	19.77	188	134735	5.00	ng	0.00
19) Chrysene-d12	23.66	240	120602	5.00	ng	0.00
25) Perylene-d12	25.34	264	95783	5.00	ng	0.00

System Monitoring Compounds						
5) Nitrobenzene-d5	10.61	82	62021	10.30	ng	0.00
9) 2-Fluorobiphenyl	14.88	172	164993	8.92	ng	0.00
21) Terphenyl-d14	22.30	244	196732	11.34	ng	0.00

Target Compounds						Ovalue
2) 1,4-Dioxane	2.74	88	262	0.06	ng	# 85
6) Naphthalene	12.29	128	2186	0.07	ng	# 93
7) 2-Methylnaphthalene	13.95	142	1177	0.06	ng	97
10) Acenaphthylene	16.07	152	6421	0.06	ng	98
11) Acenaphthene	16.49	154	1208	0.07	ng	92
12) Fluorene	17.74	166	1342	0.06	ng	100
14) Hexachlorobenzene	19.00	284	430	0.07	ng	95
15) Pentachlorophenol	19.45	266	21	Below Cal		# 62
16) Phenanthrene	19.82	178	2294	0.07	ng	99
17) Anthracene	19.91	178	1632	0.06	ng	98
18) Fluoranthene	21.72	202	1677	0.05	ng	100
20) Pyrene	22.04	202	1779	0.06	ng	99
22) Benzo(a)anthracene	23.65	228	6031	0.06	ng	# 94
23) Chrysene	23.69	228	9747	0.08	ng	94
24) Indeno(1,2,3-cd)pyrene	26.75	276	1478	0.19	ng	98
26) Benzo(b)fluoranthene	24.91	252	387	0.29	ng	# 72
27) Benzo(k)fluoranthene	24.94	252	867m	0.03	ng	
28) Benzo(a)pyrene	25.27	252	407	0.26	ng	# 66
29) Dibenzo(a,h)anthracene	26.78	278	279	0.24	ng	# 39
30) Benzo(g,h,i)perylene	27.16	276	1834	0.15	ng	# 59

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

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