

Data Path : Z:\HPCHEM1\BNA_E\Data\BE083116\
 Data File : BE092313.D
 Acq On : 31 Aug 2016 10:03
 Operator : UM/SJ
 Sample : MDL-BLK-S
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 MDL-BLK-S

Quant Time: Aug 31 14:42:34 2016
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE082516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 26 11:18:03 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	8850	0.40	ng/ul	0.00
2) Naphthalene-d8	10.49	136	45597	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.33	164	27594	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	60091	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	67037	0.40	ng/ul	0.00
20) Perylene-d12	23.63	264	64882	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.08	152	17494	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	44743	0.30	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.55	128	110	0.00	ng/ul#	1
5) 2-Methylnaphthalene	12.10	142	6	0.00	ng/ul	93
7) Acenaphthylene	14.05	152	142	0.00	ng/ul#	1
8) Acenaphthene	14.40	153	66	0.00	ng/ul#	74
9) Fluorene	15.39	166	131	0.00	ng/ul#	56
11) Pentachlorophenol	16.75	266	12	0.00	ng/ul#	6
12) Phenanthrene	17.10	178	485	0.00	ng/ul#	17
13) Anthracene	17.20	178	98	0.00	ng/ul#	1
15) Fluoranthene	19.12	202	414	0.00	ng/ul#	22
17) Pyrene	19.49	202	455	0.00	ng/ul#	100
18) Benzo(a)anthracene	21.23	228	541	0.00	ng/ul#	1
19) Chrysene	21.27	228	778	0.00	ng/ul#	1
21) Benzo(b)fluoranthene	22.90	252	218	0.00	ng/ul#	1
22) Benzo(k)fluoranthene	22.95	252	184	0.00	ng/ul#	1
23) Benzo(a)pyrene	23.51	252	182	0.00	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	26.14	276	125	0.00	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.16	278	69	0.00	ng/ul#	1
26) Benzo(g,h,i)perylene	26.91	276	43	0.00	ng/ul#	1

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

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