

Data Path : Z:\HPCHEM1\BNA_E\Data\BE060817\
 Data File : BE093074.D
 Acq On : 8 Jun 2017 10:12
 Operator : SJ/MA
 Sample : MDL-BLK-W
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 MDL-BLK-W

Quant Time: Jun 08 13:26:31 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 08 13:24:15 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	3535	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	16323	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	8400	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	24165	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	26665	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	25711	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	10095	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	26153	0.40	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.57	128	143	0.004	ng/ul#	7
5) 2-Methylnaphthalene	12.18	142	25	0.001	ng/ul#	1
7) Acenaphthylene	14.09	152	50	0.001	ng/ul#	1
8) Acenaphthene	14.42	153	63	0.002	ng/ul#	82
9) Fluorene	15.42	166	80	0.002	ng/ul#	68
11) Pentachlorophenol	16.75	266	6	0.002	ng/ul#	1
12) Phenanthrene	17.14	178	303	0.005	ng/ul#	31
13) Anthracene	17.24	178	151	0.003	ng/ul#	5
15) Fluoranthene	19.15	202	262	0.003	ng/ul#	33
17) Pyrene	19.51	202	228	0.003	ng/ul#	62
18) Benzo(a)anthracene	21.25	228	260	0.004	ng/ul#	54
19) Chrysene	21.30	228	217	0.003	ng/ul#	62
21) Benzo(b)fluoranthene	22.93	252	280	0.004	ng/ul#	1
22) Benzo(k)fluoranthene	22.98	252	268	0.003	ng/ul#	1
23) Benzo(a)pyrene	23.56	252	176	0.002	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	26.18	276	254	0.003	ng/ul#	47
25) Dibenzo(a,h)anthracene	26.20	278	206	0.003	ng/ul#	1
26) Benzo(g,h,i)perylene	26.96	276	185	0.003	ng/ul#	1

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

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