

Data Path : Z:\HPCHEM1\BNA_E\Data\BE061517\
 Data File : BE093235.D
 Acq On : 16 Jun 2017 9:57
 Operator : SJ/MA
 Sample : I3522-07DL 5X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A62DL

Quant Time: Jun 16 14:35:52 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 16 05:36:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2232	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	11206	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	6391	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	15215	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	17163	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	14597	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	601	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	1824	0.04	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.57	128	583	0.021	ng/ul#	75
5) 2-Methylnaphthalene	12.19	142	877	0.046	ng/ul	99
7) Acenaphthylene	14.09	152	35460	1.293	ng/ul	96
8) Acenaphthene	14.43	153	5986	0.271	ng/ul	98
9) Fluorene	15.42	166	15413	0.594	ng/ul	91
11) Pentachlorophenol	16.77	266	23	0.010	ng/ul#	1
12) Phenanthrene	17.14	178	514381	12.287	ng/ul#	84
13) Anthracene	17.22	178	99022	2.717	ng/ul#	86
15) Fluoranthene	19.16	202	1117057	21.759	ng/ul	84
17) Pyrene	19.52	202	859340	16.978	ng/ul#	87
18) Benzo(a)anthracene	21.24	228	529210	11.560	ng/ul#	84
19) Chrysene	21.30	228	502359	10.578	ng/ul	98
21) Benzo(b)fluoranthene	22.94	252	669541	15.115	ng/ul	87
22) Benzo(k)fluoranthene	23.12	252	58142	1.236	ng/ul	96
23) Benzo(a)pyrene	23.56	252	316026	7.545	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.19	276	166928	3.442	ng/ul#	87
25) Dibenzo(a,h)anthracene	26.20	278	55223	1.352	ng/ul	94
26) Benzo(g,h,i)perylene	26.97	276	132721	3.191	ng/ul	97

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

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