

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102617\
 Data File : BE094585.D
 Acq On : 27 Oct 2017 5:00
 Operator : SJ/JU
 Sample : I5786-01 5X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-38(FILL)_100917

Quant Time: Oct 27 10:16:18 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 25 18:44:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1547	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	9046	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.51	164	5074	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	11008	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	10977	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	11960	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	738	0.05	ng/ul	-0.02
14) Fluoranthene-d10	19.28	212	1625	0.05	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.74	128	22487	1.017	ng/ul#	93
5) 2-Methylnaphthalene	12.35	142	7881	0.509	ng/ul	98
7) Acenaphthylene	14.24	152	47137	2.054	ng/ul	96
8) Acenaphthene	14.57	153	27490	1.667	ng/ul	95
9) Fluorene	15.57	166	36591	1.976	ng/ul	95
12) Phenanthrene	17.29	178	697848	24.001	ng/ul#	88
13) Anthracene	17.38	178	173138	6.034	ng/ul#	89
15) Fluoranthene	19.32	202	1430764	42.563	ng/ul	84
17) Pyrene	19.68	202	1276411	37.614	ng/ul#	81
18) Benzo(a)anthracene	21.41	228	775650	24.718	ng/ul#	87
19) Chrysene	21.47	228	695802	21.452	ng/ul	99
21) Benzo(b)fluoranthene	23.19	252	930321m	27.653	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	341217m	9.464	ng/ul	
23) Benzo(a)pyrene	23.86	252	659131	19.600	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	26.67	276	434378	12.483	ng/ul#	75
25) Dibenzo(a,h)anthracene	26.66	278	112533	3.838	ng/ul	96
26) Benzo(g,h,i)perylene	27.51	276	320216	11.460	ng/ul	96

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

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