

Data Path : Z:\HPCHEM1\BNA_E\Data\BE032918\
 Data File : BE095909.D
 Acq On : 29 Mar 2018 11:26
 Operator : SJ/JU
 Sample : J2060-05DL 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AS8DL

Quant Time: Mar 29 17:28:57 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE031318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 29 16:32:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.42	152	1679	0.40	ng/ul	0.00
2) Naphthalene-d8	11.24	136	4933	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.00	164	3012	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.70	188	6686	0.40	ng/ul	0.00
16) Chrysene-d12	21.79	240	6769	0.40	ng/ul	0.00
20) Perylene-d12	24.42	264	7002	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.79	152	250	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.68	212	673	0.03	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	11.29	128	15975	1.160	ng/ul#	88
5) 2-Methylnaphthalene	12.87	142	13161	1.375	ng/ul	99
9) Fluorene	16.03	166	736	0.059	ng/ul#	84
12) Phenanthrene	17.75	178	25767	1.287	ng/ul#	88
13) Anthracene	17.84	178	829	0.044	ng/ul#	93
15) Fluoranthene	19.71	202	4854	0.200	ng/ul	85
17) Pyrene	20.07	202	24809	0.965	ng/ul#	79
18) Benzo(a)anthracene	21.77	228	8055	0.369	ng/ul#	88
19) Chrysene	21.83	228	18265	0.821	ng/ul	96
21) Benzo(b)fluoranthene	23.61	252	9301	0.383	ng/ul	94
22) Benzo(k)fluoranthene	23.61	252	9282	0.403	ng/ul	96
23) Benzo(a)pyrene	24.30	252	7332	0.331	ng/ul	91
24) Indeno(1,2,3-cd)pyrene	27.23	276	3758	0.153	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.24	278	2764	0.141	ng/ul#	74
26) Benzo(g,h,i)perylene	28.11	276	9630	0.457	ng/ul	93

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

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