

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE032918\
 Data File : BE095908.D
 Acq On : 29 Mar 2018 10:50
 Operator : SJ/JU
 Sample : J2060-06DL 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AS7DL

Manual Integrations
 APPROVED

Sohil
 3/29/2018 6:44:55 PM

Quant Time: Mar 29 17:17:45 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	1648	0.40	ng/ul	0.00
2) Naphthalene-d8	11.24	136	4958	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.00	164	3039	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.70	188	6669	0.40	ng/ul	0.00
16) Chrysene-d12	21.79	240	6725	0.40	ng/ul	0.00
20) Perylene-d12	24.42	264	6739	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.80	152	478	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.68	212	1317	0.06	ng/ul	0.00
Target Compounds						
					Qvalue	
3) Naphthalene	11.29	128	2278	0.165	ng/ul#	91
5) 2-Methylnaphthalene	12.87	142	986	0.102	ng/ul	98
8) Acenaphthene	15.06	153	6301	0.545	ng/ul	95
9) Fluorene	16.03	166	7661	0.611	ng/ul#	91
12) Phenanthrene	17.75	178	61947	3.102	ng/ul#	88
13) Anthracene	17.84	178	17226	0.922	ng/ul#	91
15) Fluoranthene	19.72	202	91166	3.768	ng/ul	82
17) Pyrene	20.07	202	80485	3.152	ng/ul#	78
18) Benzo(a)anthracene	21.77	228	57081	2.631	ng/ul#	85
19) Chrysene	21.83	228	41123	1.861	ng/ul	97
21) Benzo(b)fluoranthene	23.61	252	46751m	2.003	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	15797m	0.712	ng/ul	
23) Benzo(a)pyrene	24.30	252	32222	1.512	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	27.23	276	18102	0.766	ng/ul#	74
25) Dibenzo(a,h)anthracene	27.24	278	5579	0.295	ng/ul	95
26) Benzo(g,h,i)perylene	28.11	276	15778	0.777	ng/ul	94

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

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