

Data Path : Z:\HPCHEM1\BNA E\DATA\BE031518\
 Data File : BE095808.D
 Acq On : 15 Mar 2018 17:52
 Operator : SJ/JU
 Sample : J1835-12DL 10X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6F08DL

Manual Integrations
 APPROVED

Sohil
 3/16/2018 12:03:17 PM

Quant Time: Mar 16 03:07:37 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE031318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 16 02:40:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.45	152	1895	0.40	ng/ul	0.00
2) Naphthalene-d8	11.28	136	5450	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.04	164	2962	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.74	188	6455m	0.40	ng/ul	0.00
16) Chrysene-d12	21.84	240	6281m	0.40	ng/ul	0.00
20) Perylene-d12	24.49	264	6281	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.83	152	211	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.72	212	533m	0.03	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.33	128	797	0.052	ng/ul#	75
5) 2-Methylnaphthalene	12.90	142	1271	0.120	ng/ul	100
12) Phenanthrene	17.78	178	8382m	0.434	ng/ul	
13) Anthracene	17.87	178	504m	0.028	ng/ul	
15) Fluoranthene	19.75	202	2194m	0.094	ng/ul	
17) Pyrene	20.10	202	6838	0.287	ng/ul#	80
18) Benzo(a)anthracene	21.82	228	2749m	0.136	ng/ul	
19) Chrysene	21.87	228	5275m	0.256	ng/ul	
21) Benzo(b)fluoranthene	23.67	252	2128m	0.098	ng/ul	
22) Benzo(k)fluoranthene	23.71	252	500m	0.024	ng/ul	
23) Benzo(a)pyrene	24.38	252	1958	0.099	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	27.34	276	1087	0.049	ng/ul#	59
26) Benzo(g,h,i)perylene	28.24	276	2376	0.126	ng/ul#	23

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

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