

Data Path : Z:\HPCHEM1\BNA E\DATA\BE031518\
 Data File : BE095812.D
 Acq On : 15 Mar 2018 20:16
 Operator : SJ/JU
 Sample : J1813-04DL 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6F20DL

Manual Integrations
 APPROVED

Sohil
 3/16/2018 12:03:28 PM

Quant Time: Mar 16 04:00:52 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	1980	0.40	ng/ul	0.00
2) Naphthalene-d8	11.27	136	5869	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.03	164	3188	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.74	188	6601	0.40	ng/ul	0.00
16) Chrysene-d12	21.84	240	7056	0.40	ng/ul	0.00
20) Perylene-d12	24.49	264	7094	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.83	152	753	0.08	ng/ul	0.00
14) Fluoranthene-d10	19.72	212	1764	0.08	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.32	128	2886	0.176	ng/ul#	90
5) 2-Methylnaphthalene	12.90	142	903	0.079	ng/ul	98
7) Acenaphthylene	14.76	152	1547	0.096	ng/ul#	90
8) Acenaphthene	15.09	153	3486	0.288	ng/ul	93
9) Fluorene	16.06	166	2298	0.175	ng/ul	87
12) Phenanthrene	17.78	178	34279	1.734	ng/ul#	89
13) Anthracene	17.87	178	7699	0.417	ng/ul#	93
15) Fluoranthene	19.75	202	75548	3.155	ng/ul	81
17) Pyrene	20.10	202	63343m	2.364	ng/ul	
18) Benzo(a)anthracene	21.81	228	49105	2.157	ng/ul#	88
19) Chrysene	21.87	228	41689	1.798	ng/ul	94
21) Benzo(b)fluoranthene	23.67	252	64274m	2.616	ng/ul	
22) Benzo(k)fluoranthene	23.71	252	18661m	0.799	ng/ul	
23) Benzo(a)pyrene	24.37	252	40862	1.821	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	27.35	276	27883	1.121	ng/ul#	68
25) Dibenzo(a,h)anthracene	27.35	278	6391	0.321	ng/ul#	65
26) Benzo(a,h,i)perylene	28.25	276	23840	1.116	ng/ul#	92

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

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