

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042018\
 Data File : BE096064.D
 Acq On : 21 Apr 2018 5:25
 Operator : SJ/JU
 Sample : J2434-03
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB8

Manual Integrations
 APPROVED

Sohil
 4/23/2018 6:44:20 PM

Quant Time: Apr 23 02:28:10 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 23 01:07:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	2005	0.40	ng/ul	0.00
2) Naphthalene-d8	11.20	136	5542	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	3307	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	7388m	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	7330m	0.40	ng/ul	0.00
20) Perylene-d12	24.38	264	7545	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.76	152	1640	0.16	ng/ul	0.00
14) Fluoranthene-d10	19.66	212	4672m	0.15	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.25	128	438	0.028	ng/ul#	61
7) Acenaphthylene	14.70	152	376	0.022	ng/ul#	64
8) Acenaphthene	15.03	153	361	0.029	ng/ul#	92
9) Fluorene	16.00	166	438	0.032	ng/ul#	84
12) Phenanthrene	17.72	178	11047m	0.507	ng/ul	
13) Anthracene	17.80	178	2104m	0.097	ng/ul	
15) Fluoranthene	19.68	202	34747m	1.173	ng/ul	
17) Pyrene	20.04	202	31798	1.277	ng/ul#	78
18) Benzo(a)anthracene	21.75	228	21184m	0.715	ng/ul	
19) Chrysene	21.80	228	16513m	0.775	ng/ul	
21) Benzo(b)fluoranthene	23.57	252	18929m	0.775	ng/ul	
22) Benzo(k)fluoranthene	23.62	252	7239m	0.298	ng/ul	
23) Benzo(a)pyrene	24.26	252	14236	0.604	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.16	276	8428	0.322	ng/ul#	74
25) Dibenzo(a,h)anthracene	27.17	278	2237	0.106	ng/ul#	87
26) Benzo(a,h,i)perylene	28.03	276	7573	0.346	ng/ul	95

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

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