

Data Path : Z:\HPCHEM1\BNA E\DATA\BE050318\
 Data File : BE096334.D
 Acq On : 3 May 2018 21:53
 Operator : SJ/JU
 Sample : J2440-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Y18

Manual Integrations
 APPROVED

Sohil
 5/4/2018 1:35:06 PM

Quant Time: May 04 03:47:08 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE050318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 04 01:53:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.35	152	638	0.40	ng/ul	0.00
2) Naphthalene-d8	11.17	136	2776	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	1806	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.65	188	4642m	0.40	ng/ul	0.00
16) Chrysene-d12	21.74	240	5012m	0.40	ng/ul	0.00
20) Perylene-d12	24.34	264	4115m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	1134	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.63	212	3833m	0.19	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.67	152	501	0.053	ng/ul#	77
9) Fluorene	15.97	166	186	0.020	ng/ul#	81
12) Phenanthrene	17.69	178	5077m	0.363	ng/ul	
13) Anthracene	17.78	178	503m	0.037	ng/ul	
15) Fluoranthene	19.66	202	12185m	0.613	ng/ul	
17) Pyrene	20.02	202	3159m	0.560	ng/ul	
18) Benzo(a)anthracene	21.73	228	1767m	0.226	ng/ul	
19) Chrysene	21.78	228	5094m	0.340	ng/ul	
21) Benzo(b)fluoranthene	23.54	252	5833m	0.414	ng/ul	
22) Benzo(k)fluoranthene	23.59	252	1854m	0.136	ng/ul	
23) Benzo(a)pyrene	24.23	252	3263m	0.247	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.12	276	2404m	0.165	ng/ul	
25) Dibenzo(a,h)anthracene	27.12	278	618m	0.052	ng/ul	
26) Benzo(g,h,i)perylene	27.98	276	2145m	0.174	ng/ul	

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

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