

Data Path : Z:\HPCHEM1\BNA E\DATA\BE032818\
 Data File : BE095874.D
 Acq On : 28 Mar 2018 13:34
 Operator : SJ/JU
 Sample : J2060-08
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AT5

Manual Integrations
 APPROVED

Sohil
 3/29/2018 5:17:37 PM

Quant Time: Mar 28 18:03:20 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE031318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 28 16:31:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.42	152	1647	0.40	ng/ul	0.00
2) Naphthalene-d8	11.24	136	4670	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.00	164	3220	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.70	188	7427m	0.40	ng/ul	0.00
16) Chrysene-d12	21.79	240	7907m	0.40	ng/ul	0.00
20) Perylene-d12	24.42	264	7934m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.80	152	3306	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.68	212	10921m	0.46	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.29	128	1269	0.097	ng/ul#	86
5) 2-Methylnaphthalene	12.87	142	491	0.054	ng/ul	98
8) Acenaphthene	15.06	153	2983	0.244	ng/ul	93
9) Fluorene	16.02	166	3552	0.267	ng/ul	89
12) Phenanthrene	17.75	178	23411m	1.053	ng/ul	
13) Anthracene	17.83	178	7548m	0.363	ng/ul	
15) Fluoranthene	19.71	202	37985m	1.410	ng/ul	
17) Pyrene	20.07	202	36178	1.205	ng/ul#	77
18) Benzo(a)anthracene	21.77	228	21761	0.853	ng/ul#	87
19) Chrysene	21.83	228	17216	0.663	ng/ul	98
21) Benzo(b)fluoranthene	23.60	252	18589m	0.676	ng/ul	
22) Benzo(k)fluoranthene	23.65	252	6204m	0.237	ng/ul	
23) Benzo(a)pyrene	24.30	252	13579m	0.541	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.21	276	7197m	0.259	ng/ul	
25) Dibenzo(a,h)anthracene	27.21	278	2283m	0.103	ng/ul	
26) Benzo(a,h,i)perylene	28.09	276	6464m	0.270	ng/ul	

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

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