

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042318\
 Data File : BE096082.D
 Acq On : 23 Apr 2018 13:23
 Operator : SJ/JU
 Sample : J2355-06MS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Y33MS

Manual Integrations
 APPROVED

Sohil
 4/24/2018 7:09:50 PM

Quant Time: Apr 24 10:32:03 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 24 05:21:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	1883	0.40	ng/ul	0.00
2) Naphthalene-d8	11.20	136	5213	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	2892	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	6277	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	7773	0.40	ng/ul	0.00
20) Perylene-d12	24.38	264	6315	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.76	152	3184	0.33	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	8331	0.32	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.25	128	7540	0.518	ng/ul#	89
5) 2-Methylnaphthalene	12.83	142	1815	0.184	ng/ul	100
7) Acenaphthylene	14.70	152	2882	0.192	ng/ul	97
8) Acenaphthene	15.03	153	6313	0.587	ng/ul	96
9) Fluorene	15.99	166	2790	0.230	ng/ul	91
11) Pentachlorophenol	17.33	266	368	0.277	ng/ul	98
12) Phenanthrene	17.71	178	39380	2.129	ng/ul#	88
13) Anthracene	17.80	178	6249	0.339	ng/ul#	91
15) Fluoranthene	19.69	202	72834	2.893	ng/ul	82
17) Pyrene	20.04	202	71884	2.722	ng/ul#	77
18) Benzo(a)anthracene	21.75	228	40415m	1.287	ng/ul	
19) Chrysene	21.80	228	32706m	1.447	ng/ul	
21) Benzo(b)fluoranthene	23.57	252	42866m	2.098	ng/ul	
22) Benzo(k)fluoranthene	23.62	252	12088m	0.594	ng/ul	
23) Benzo(a)pyrene	24.26	252	28420	1.441	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.17	276	20121	0.918	ng/ul#	73
25) Dibenzo(a,h)anthracene	27.17	278	4930	0.279	ng/ul	96
26) Benzo(g,h,i)perylene	28.04	276	15560	0.850	ng/ul	94

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

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