

Data Path : Z:\HPCHEM1\BNA E\DATA\BE012818\
 Data File : BE095228.D
 Acq On : 29 Jan 2018 3:51
 Operator : SJ/JU
 Sample : J1223-07
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6A73

Manual Integrations
 APPROVED

Sohil
 1/29/2018 6:26:08 PM

Quant Time: Jan 29 08:35:36 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 29 08:11:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2678	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	8281	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	5276	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	10976	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	12324	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	11224m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	3100	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	9486	0.23	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.32	128	42943	1.852	ng/ul#	89
5) 2-Methylnaphthalene	12.89	142	6318	0.408	ng/ul	100
7) Acenaphthylene	14.75	152	1181	0.047	ng/ul#	85
8) Acenaphthene	15.08	153	4304	0.219	ng/ul	94
9) Fluorene	16.03	166	2513	0.124	ng/ul#	76
11) Pentachlorophenol	17.36	266	62097	29.210	ng/ul	95
12) Phenanthrene	17.75	178	75776	2.053	ng/ul#	86
13) Anthracene	17.85	178	10847	0.321	ng/ul#	91
15) Fluoranthene	19.73	202	188820	3.950	ng/ul	83
17) Pyrene	20.08	202	139843	3.382	ng/ul#	79
18) Benzo(a)anthracene	21.79	228	34944	0.907	ng/ul#	85
19) Chrysene	21.84	228	35619	0.890	ng/ul	97
21) Benzo(b)fluoranthene	23.62	252	27778m	0.701	ng/ul	
22) Benzo(k)fluoranthene	23.67	252	7087m	0.196	ng/ul	
23) Benzo(a)pyrene	24.32	252	9868m	0.279	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.24	276	4950m	0.128	ng/ul	
25) Dibenzo(a,h)anthracene	27.26	278	1257m	0.040	ng/ul	
26) Benzo(g,h,i)perylene	28.12	276	3623m	0.110	ng/ul	

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

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