

Data Path : Z:\HPCHEM1\BNA E\DATA\BE012818\
 Data File : BE095223.D
 Acq On : 29 Jan 2018 00:52
 Operator : SJ/JU
 Sample : J1223-02
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6A68

Manual Integrations
 APPROVED

Sohil
 1/29/2018 6:26:01 PM

Quant Time: Jan 29 08:18:17 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	2874	0.40	ng/ul	0.00
2) Naphthalene-d8	11.27	136	9423	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	6654	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.73	188	12672m	0.40	ng/ul	0.02
16) Chrysene-d12	21.80	240	12875	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	11930m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.82	152	3855	0.25	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	11144	0.23	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.31	128	1111	0.042	ng/ul#	57
5) 2-Methylnaphthalene	12.89	142	46607	2.645	ng/ul	100
7) Acenaphthylene	14.74	152	5420	0.169	ng/ul#	33
8) Acenaphthene	15.08	153	43712	1.760	ng/ul	94
9) Fluorene	16.02	166	25289	0.987	ng/ul#	65
11) Pentachlorophenol	17.36	266	94207	38.383	ng/ul	96
12) Phenanthrene	17.75	178	143992	3.379	ng/ul#	86
13) Anthracene	17.84	178	18321	0.470	ng/ul	93
15) Fluoranthene	19.73	202	212067	3.843	ng/ul	82
17) Pyrene	20.08	202	187697	4.345	ng/ul#	79
18) Benzo(a)anthracene	21.79	228	46369	1.152	ng/ul#	87
19) Chrysene	21.84	228	49889	1.194	ng/ul	94
21) Benzo(b)fluoranthene	23.62	252	57642m	1.370	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	15134m	0.393	ng/ul	
23) Benzo(a)pyrene	24.32	252	16295m	0.434	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.25	276	12379m	0.300	ng/ul	
25) Dibenzo(a,h)anthracene	27.27	278	2863m	0.085	ng/ul	
26) Benzo(g,h,i)perylene	28.12	276	9693m	0.278	ng/ul	

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

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