

Data Path : Z:\HPCHEM1\BNA E\DATA\BE012818\
 Data File : BE095235.D
 Acq On : 29 Jan 2018 8:02
 Operator : SJ/JU
 Sample : J1223-16
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6B02

Manual Integrations
 APPROVED

Sohil
 1/29/2018 6:26:28 PM

Quant Time: Jan 29 13:32:09 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	3607	0.40	ng/ul	0.00
2) Naphthalene-d8	11.27	136	11675	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	9025	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.73	188	15165m	0.40	ng/ul	0.02
16) Chrysene-d12	21.81	240	15212	0.40	ng/ul	0.00
20) Perylene-d12	24.45	264	13505m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.82	152	6929	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	17958	0.31	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.32	128	57919	1.772	ng/ul#	88
5) 2-Methylnaphthalene	12.89	142	59780	2.738	ng/ul	100
7) Acenaphthylene	14.75	152	3715	0.086	ng/ul#	75
8) Acenaphthene	15.08	153	63716	1.891	ng/ul	97
9) Fluorene	16.03	166	72538	2.088	ng/ul#	68
11) Pentachlorophenol	17.36	266	19410	6.608	ng/ul	95
12) Phenanthrene	17.75	178	406274	7.966	ng/ul#	84
13) Anthracene	17.85	178	42643	0.913	ng/ul#	90
15) Fluoranthene	19.73	202	268357	4.064	ng/ul	83
17) Pyrene	20.09	202	175057	3.429	ng/ul#	76
18) Benzo(a)anthracene	21.79	228	39176	0.824	ng/ul#	85
19) Chrysene	21.84	228	28410	0.575	ng/ul	96
21) Benzo(b)fluoranthene	23.63	252	19401m	0.407	ng/ul	
22) Benzo(k)fluoranthene	23.67	252	5704m	0.131	ng/ul	
23) Benzo(a)pyrene	24.33	252	5838m	0.137	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.25	276	3393m	0.073	ng/ul	
25) Dibenzo(a,h)anthracene	27.26	278	1048m	0.027	ng/ul	
26) Benzo(g,h,i)perylene	28.12	276	2976m	0.075	ng/ul	

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

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