

Data Path : U:\HPCHEM1\BNA E\DATA\BE020218\
 Data File : BE095404.D
 Acq On : 2 Feb 2018 19:24
 Operator : SJ/JU
 Sample : J1244-02DL 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6B43DL

Manual Integrations
 APPROVED

Sohil
 2/5/2018 5:48:08 PM

Quant Time: Feb 03 05:15:26 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 03 04:31:32 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	1380m	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	5720	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	3658	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	7514	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7870	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	7778	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	493	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	1245	0.04	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.31	128	97945	6.115	ng/ul#	88
5) 2-Methylnaphthalene	12.89	142	53531	5.005	ng/ul	100
7) Acenaphthylene	14.74	152	939	0.053	ng/ul#	66
8) Acenaphthene	15.08	153	42615	3.121	ng/ul	95
9) Fluorene	16.02	166	35775	2.541	ng/ul#	82
11) Pentachlorophenol	17.36	266	11204	7.698	ng/ul	97
12) Phenanthrene	17.75	178	166326	6.582	ng/ul#	88
13) Anthracene	17.85	178	22427	0.970	ng/ul#	92
15) Fluoranthene	19.72	202	96592	2.952	ng/ul	82
17) Pyrene	20.08	202	79675	3.017	ng/ul#	76
18) Benzo(a)anthracene	21.79	228	14572	0.592	ng/ul#	87
19) Chrysene	21.84	228	10885	0.426	ng/ul	97
21) Benzo(b)fluoranthene	23.62	252	7454m	0.272	ng/ul	
22) Benzo(k)fluoranthene	23.67	252	2636m	0.105	ng/ul	
23) Benzo(a)pyrene	24.32	252	3724m	0.152	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.24	276	1263	0.047	ng/ul#	77
26) Benzo(a,h,i)perylene	28.12	276	1059	0.047	ng/ul#	80

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

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