

Data Path : Z:\HPCHEM1\BNA E\DATA\BE013018\
 Data File : BE095307.D
 Acq On : 31 Jan 2018 9:00
 Operator : SJ/JU
 Sample : J1233-12DL 10X
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6B16DL

Manual Integrations
 APPROVED

Sohil
 1/31/2018 8:28:54 PM

Quant Time: Jan 31 10:00:54 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 31 08:16:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2991	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	9651	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	6241	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	12630	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	13216	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	12582	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	260	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	646	0.01	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.31	128	38718	1.433	ng/ul#	88
5) 2-Methylnaphthalene	12.89	142	39620	2.196	ng/ul	100
7) Acenaphthylene	14.74	152	897	0.030	ng/ul#	41
8) Acenaphthene	15.08	153	19301	0.829	ng/ul	94
9) Fluorene	16.03	166	26514	1.104	ng/ul#	80
11) Pentachlorophenol	17.36	266	217	0.089	ng/ul	92
12) Phenanthrene	17.75	178	162947	3.836	ng/ul#	88
13) Anthracene	17.85	178	24473	0.629	ng/ul#	92
15) Fluoranthene	19.72	202	102498	1.864	ng/ul	84
17) Pyrene	20.08	202	79579	1.794	ng/ul#	77
18) Benzo(a)anthracene	21.79	228	19237	0.466	ng/ul#	86
19) Chrysene	21.84	228	11940	0.278	ng/ul	96
21) Benzo(b)fluoranthene	23.62	252	7645m	0.172	ng/ul	
22) Benzo(k)fluoranthene	23.67	252	3148m	0.078	ng/ul	
23) Benzo(a)pyrene	24.32	252	4577	0.115	ng/ul	89
24) Indeno(1,2,3-cd)pyrene	27.23	276	1419	0.033	ng/ul#	74

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

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