

Data Path : U:\HPCHEM1\BNA E\DATA\BE012618\
 Data File : BE095157.D
 Acq On : 27 Jan 2018 8:39
 Operator : SJ/JU
 Sample : J1174-02
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6A01

Manual Integrations
 APPROVED

Sohil
 1/29/2018 6:34:35 PM

Quant Time: Jan 29 09:25:41 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 Jan 27 03:18:43 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	3584	0.40	ng/ul	0.00
2) Naphthalene-d8	11.27	136	11352	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	7141	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.73	188	16312m	0.40	ng/ul	0.02
16) Chrysene-d12	21.80	240	17627	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	16313m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.82	152	3845	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	13063	0.21	ng/ul	0.00
Target Compounds				Ovalue		
11) Pentachlorophenol	17.36	266	2827	0.895	ng/ul	97
12) Phenanthrene	17.75	178	1312	0.024	ng/ul#	87
13) Anthracene	17.84	178	1385	0.028	ng/ul#	87
15) Fluoranthene	19.73	202	6172	0.087	ng/ul	84
17) Pyrene	20.08	202	7400m	0.125	ng/ul	
18) Benzo(a)anthracene	21.79	228	2177	0.040	ng/ul#	78
19) Chrysene	21.84	228	5949	0.104	ng/ul#	86
21) Benzo(b)fluoranthene	23.62	252	6342m	0.110	ng/ul	
22) Benzo(k)fluoranthene	23.67	252	1343m	0.026	ng/ul	

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

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