

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE071818\
 Data File : BE096995.D
 Acq On : 18 Jul 2018 16:02
 Operator : SJ/JU
 Sample : J3967-19
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 DB1F9

Manual Integrations
 APPROVED

Sohil
 7/19/2018 11:36:53 AM

Quant Time: Jul 18 16:51:31 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE071718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 18 15:28:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1636	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	9030	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	6260	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	15463m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	17560m	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	14339m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	4905	0.34	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	15920m	0.34	ng/ul	0.00
Target Compounds				Qvalue		
8) Acenaphthene	14.08	153	3057	0.144	ng/ul	95
9) Fluorene	15.08	166	7580	0.337	ng/ul	93
11) Pentachlorophenol	16.44	266	541m	0.310	ng/ul	
12) Phenanthrene	16.82	178	24720m	0.609	ng/ul	
13) Anthracene	16.91	178	10026m	0.283	ng/ul	
15) Fluoranthene	18.84	202	39040m	0.729	ng/ul	
17) Pyrene	19.21	202	27266	0.537	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	2606m	0.056	ng/ul	
19) Chrysene	21.01	228	7647	0.145	ng/ul#	92
21) Benzo(b)fluoranthene	22.54	252	5610m	0.138	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	1678m	0.040	ng/ul	
23) Benzo(a)pyrene	23.12	252	1705	0.048	ng/ul	91
24) Indeno(1,2,3-cd)pyrene	25.50	276	2051	0.051	ng/ul#	81
26) Benzo(g,h,i)perylene	26.20	276	2189	0.061	ng/ul#	92

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

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