

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE070618\
 Data File : BE096864.D
 Acq On : 6 Jul 2018 21:48
 Operator : SJ/JU
 Sample : J3721-08
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F1H19

Manual Integrations
 APPROVED

Sohil
 7/9/2018 4:14:10 PM

Quant Time: Jul 09 05:31:22 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE070618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 06 16:55:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	4186	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	22240	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	14061	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	33631m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	32847m	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	28931m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	9961	0.27	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	25643m	0.31	ng/ul	0.00
Target Compounds						Qvalue
7) Acenaphthylene	13.74	152	6702	0.116	ng/ul	98
8) Acenaphthene	14.09	153	1600	0.034	ng/ul	94
9) Fluorene	15.08	166	1991	0.041	ng/ul	93
12) Phenanthrene	16.82	178	10791m	0.124	ng/ul	
13) Anthracene	16.91	178	15504m	0.211	ng/ul	
15) Fluoranthene	18.84	202	69925m	0.757	ng/ul	
17) Pyrene	19.21	202	120549	1.118	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	85833m	1.108	ng/ul	
19) Chrysene	21.01	228	133715	1.332	ng/ul	96
21) Benzo(b)fluoranthene	22.54	252	191984m	2.431	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	60854m	0.706	ng/ul	
23) Benzo(a)pyrene	23.12	252	105474m	1.579	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.50	276	58865m	0.849	ng/ul	
25) Dibenzo(a,h)anthracene	25.51	278	14177m	0.257	ng/ul	
26) Benzo(g,h,i)perylene	26.20	276	48010m	0.713	ng/ul	

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

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