

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082818\
 Data File : BF108561.D
 Acq On : 28 Aug 2018 17:51
 Operator : JU/SJ
 Sample : J4276-05 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-02-082718

Manual Integrations
 APPROVED

Sohil
 8/29/2018 3:11:31 PM

Quant Time: Aug 29 10:38:16 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 22 11:01:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.00	152	111636	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	434293	20.00	ng	0.00
38) Acenaphthene-d10	10.04	164	200142	20.00	ng	-0.01
63) Phenanthrene-d10	11.54	188	333321	20.00	ng	0.00
75) Chrysene-d12	14.19	240	279463	20.00	ng	0.00
86) Perylene-d12	15.75	264	231862	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	37065	6.01	ng	0.03
7) Phenol-d6	6.64	99	61145	7.46	ng	0.01
23) Nitrobenzene-d5	7.57	82	38588	4.96	ng	0.00
41) 2,4,6-Tribromophenol	10.84	330	12147	6.08	ng	-0.01
44) 2-Fluorobiphenyl	9.35	172	84118	6.75	ng	-0.01
78) Terphenyl-d14	13.12	244	67350	6.13	ng	0.00
Target Compounds						
48) Acenaphthylene	9.91	152	40504	2.197	ng	98
57) Fluorene	10.59	166	54164	4.182	ng	99
70) Phenanthrene	11.57	178	283777	18.050	ng	99
71) Anthracene	11.62	178	82480	5.119	ng	97
74) Fluoranthene	12.76	202	322147	18.775	ng	96
77) Pyrene	12.99	202	329637	18.631	ng	99
80) Benzo(a)anthracene	14.18	228	154774	9.650	ng	98
82) Chrysene	14.22	228	146691	9.976	ng	97
85) Indeno(1,2,3-cd)pyrene	17.37	276	54562	4.283	ng	# 87
87) Benzo(b)fluoranthene	15.29	252	157549m	11.372	ng	
88) Benzo(k)fluoranthene	15.31	252	48320m	3.794	ng	
89) Benzo(a)pyrene	15.69	252	117175m	9.445	ng	
91) Benzo(a,h,i)perylene	17.87	276	56473	5.587	ng	# 89

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

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