

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082218\
 Data File : BF108397.D
 Acq On : 22 Aug 2018 21:39
 Operator : JU/SJ
 Sample : J4276-03 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-01-082118

Manual Integrations
 APPROVED

Sohil
 8/23/2018 3:43:23 PM

Quant Time: Aug 23 12:04:52 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	143141	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	553974	20.00	ng	0.00
38) Acenaphthene-d10	10.05	164	259149	20.00	ng	0.00
63) Phenanthrene-d10	11.54	188	423787	20.00	ng	0.00
75) Chrysene-d12	14.19	240	298180	20.00	ng	0.00
86) Perylene-d12	15.75	264	272151	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	153841	19.46	ng	-0.01
7) Phenol-d6	6.61	99	215257	20.49	ng	-0.02
23) Nitrobenzene-d5	7.56	82	131673	13.26	ng	-0.01
41) 2,4,6-Tribromophenol	10.84	330	39853	15.42	ng	-0.01
44) 2-Fluorobiphenyl	9.35	172	277080	17.17	ng	-0.01
78) Terphenyl-d14	13.12	244	195479	16.67	ng	0.00
Target Compounds						
57) Fluorene	10.60	166	37166	2.216	ng	Qvalue 97
70) Phenanthrene	11.57	178	247345	12.374	ng	99
71) Anthracene	11.62	178	63095	3.080	ng	96
74) Fluoranthene	12.76	202	286148	13.117	ng	95
77) Pyrene	13.00	202	265247	14.051	ng	98
80) Benzo(a)anthracene	14.18	228	117608	6.873	ng	98
82) Chrysene	14.22	228	106242	6.772	ng	97
85) Indeno(1,2,3-cd)pyrene	17.36	276	39010	2.870	ng	# 82
87) Benzo(b)fluoranthene	15.28	252	112496m	6.918	ng	
88) Benzo(k)fluoranthene	15.31	252	39113m	2.616	ng	
89) Benzo(a)pyrene	15.68	252	83229	5.715	ng	96
91) Benzo(a,h,i)perylene	17.85	276	38784	3.269	ng	# 89

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

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