

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\
 Data File : BF110735.D
 Acq On : 13 Nov 2018 22:38
 Operator : JU/SJ
 Sample : J5770-01 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-01-111218-A

Manual Integrations
 APPROVED

Sohil
 11/14/2018 1:52:29 PM

Quant Time: Nov 14 06:25:12 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	83441	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	314824	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	124430	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	199039	20.00	ng	0.00
76) Chrysene-d12	14.13	240	171545	20.00	ng	-0.01
87) Perylene-d12	15.66	264	115898	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	277923	54.10	ng	0.00
7) Phenol-d6	6.57	99	342695	52.17	ng	0.00
23) Nitrobenzene-d5	7.51	82	204383	37.39	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	55473	44.24	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	338796	38.89	ng	0.00
79) Terphenyl-d14	13.06	244	260588	28.45	ng	-0.01
Target Compounds						
10) Phenol	6.58	94	15833	2.079	ng	Qvalue # 72
71) Phenanthrene	11.50	178	33396	3.132	ng	98
75) Fluoranthene	12.70	202	66364	6.208	ng	98
78) Pyrene	12.93	202	68423	5.180	ng	97
81) Benzo(a)anthracene	14.12	228	34296	3.345	ng	97
83) Chrysene	14.16	228	36291	3.715	ng	97
86) Indeno(1,2,3-cd)pyrene	17.23	276	14110	2.013	ng	# 92
88) Benzo(b)fluoranthene	15.20	252	32817m	4.733	ng	
90) Benzo(a)pyrene	15.59	252	24059	3.819	ng	# 93
92) Benzo(g,h,i)perylene	17.72	276	13135	2.483	ng	# 87

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

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