

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\
 Data File : BF110736.D
 Acq On : 13 Nov 2018 23:06
 Operator : JU/SJ
 Sample : J5770-03 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

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

Quant Time: Nov 14 06:28:12 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	86289	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	331264	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	126761	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	207350	20.00	ng	0.00
76) Chrysene-d12	14.13	240	175799	20.00	ng	0.00
87) Perylene-d12	15.66	264	118129	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	304916	57.40	ng	0.00
7) Phenol-d6	6.57	99	376451	55.42	ng	0.00
23) Nitrobenzene-d5	7.51	82	226233	39.33	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	62116	48.63	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	358419	40.38	ng	0.00
79) Terphenyl-d14	13.06	244	287212	30.60	ng	-0.01
Target Compounds						
10) Phenol	6.58	94	17951	2.279	ng	Qvalue # 73
75) Fluoranthene	12.70	202	41139	3.694	ng	99
78) Pyrene	12.93	202	47544	3.512	ng	98
81) Benzo(a)anthracene	14.12	228	21027	2.001	ng	91
83) Chrysene	14.16	228	24125	2.410	ng	98
88) Benzo(b)fluoranthene	15.20	252	25140m	3.557	ng	
90) Benzo(a)pyrene	15.59	252	16429	2.559	ng	# 91
92) Benzo(g,h,i)perylene	17.72	276	12560	2.330	ng	# 81

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

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