

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111418\
 Data File : BF110765.D
 Acq On : 14 Nov 2018 15:39
 Operator : JU/SJ
 Sample : J5930-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 32727-45914-21716

Manual Integrations
 APPROVED

Sohil
 11/15/2018 10:59:24 AM

Quant Time: Nov 14 16:25:57 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	85897	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	374985	20.00	ng	0.01
39) Acenaphthene-d10	9.99	164	158980	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	284902	20.00	ng	0.00
76) Chrysene-d12	14.13	240	176276	20.00	ng	-0.01
87) Perylene-d12	15.66	264	166310	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	566012	107.03	ng	0.00
7) Phenol-d6	6.57	99	773755	114.42	ng	0.00
23) Nitrobenzene-d5	7.54	82	470268	72.22	ng	0.02
42) 2,4,6-Tribromophenol	10.78	330	160615	100.26	ng	0.00
45) 2-Fluorobiphenyl	9.32	172	735730	66.09	ng	0.01
79) Terphenyl-d14	13.06	244	592981	63.00	ng	0.00
Target Compounds						
6) Aniline	6.61	93	522368	59.234	ng	Qvalue # 55
10) Phenol	6.59	94	22477	2.867	ng	94
31) Naphthalene	8.26	128	81536m	4.632	ng	
50) Dimethylphthalate	9.69	163	74921	6.316	ng	# 99
71) Phenanthrene	11.50	178	52543	3.442	ng	100
75) Fluoranthene	12.70	202	66996	4.378	ng	99
77) Benzydine	12.82	184	14359	2.962	ng	97
78) Pyrene	12.93	202	56733	4.180	ng	97
81) Benzo(a)anthracene	14.12	228	27244	2.586	ng	96
83) Chrysene	14.16	228	26951	2.685	ng	98
88) Benzo(b)fluoranthene	15.20	252	29615m	2.977	ng	
90) Benzo(a)pyrene	15.59	252	23155	2.562	ng	# 42

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

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