

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121418\
 Data File : BF111427.D
 Acq On : 14 Dec 2018 19:07
 Operator : JU/SJ
 Sample : J6376-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEST-PIT-2

Manual Integrations
 APPROVED

Sohil
 12/17/2018 2:36:39 PM

Quant Time: Dec 14 22:21:32 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 14 13:26:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	161392	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	636467	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	288577	20.00	ng	0.00
64) Phenanthrene-d10	11.31	188	411529	20.00	ng	0.00
76) Chrysene-d12	13.95	240	313674	20.00	ng	0.00
87) Perylene-d12	15.39	264	251317	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1097285	113.14	ng	0.00
7) Phenol-d6	6.44	99	1416881	109.83	ng	0.00
23) Nitrobenzene-d5	7.35	82	855519	80.04	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	245651	95.03	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1482071	79.33	ng	0.00
79) Terphenyl-d14	12.90	244	943774	70.65	ng	0.00
Target Compounds						
10) Phenol	6.45	94	93467	6.504	ng	88
50) Dimethylphthalate	9.55	163	138072	6.644	ng	99
75) Fluoranthene	12.52	202	194477	9.375	ng	96
78) Pyrene	12.75	202	339372	15.346	ng	95
81) Benzo(a)anthracene	13.94	228	153002m	8.971	ng	
83) Chrysene	13.97	228	157174	8.748	ng	97
86) Indeno(1,2,3-cd)pyrene	16.80	276	34050	2.626	ng	# 88
88) Benzo(b)fluoranthene	14.97	252	113999m	7.780	ng	
89) Benzo(k)fluoranthene	15.00	252	35529m	2.538	ng	
90) Benzo(a)pyrene	15.33	252	92164	6.975	ng	99
92) Benzo(g,h,i)perylene	17.22	276	38274	3.741	ng	93

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

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