

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
 Data File : BF111474.D
 Acq On : 15 Dec 2018 17:48
 Operator : JU/SJ
 Sample : J6361-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TN4803-121218

Manual Integrations
 APPROVED

Sohil
 12/17/2018 3:12:14 PM

Quant Time: Dec 16 01:39:57 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	165667	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	648111	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	273542	20.00	ng	0.00
64) Phenanthrene-d10	11.31	188	457749	20.00	ng	0.00
76) Chrysene-d12	13.95	240	418293	20.00	ng	0.00
87) Perylene-d12	15.39	264	313441	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1002753	100.72	ng	0.00
7) Phenol-d6	6.44	99	1271913	96.05	ng	0.00
23) Nitrobenzene-d5	7.35	82	770405	70.79	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	196228	80.08	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1282858	72.44	ng	0.00
79) Terphenyl-d14	12.90	244	1076406	60.43	ng	0.00
Target Compounds						
10) Phenol	6.45	94	88307	5.986	ng	94
50) Dimethylphthalate	9.55	163	96591	4.903	ng	97
71) Phenanthrene	11.33	178	128504	5.448	ng	# 92
75) Fluoranthene	12.52	202	252666	10.950	ng	97
78) Pyrene	12.75	202	226717	7.688	ng	95
81) Benzo(a)anthracene	13.94	228	97576	4.290	ng	100
83) Chrysene	13.97	228	112986	4.716	ng	96
86) Indeno(1,2,3-cd)pyrene	16.80	276	37785	2.185	ng	96
88) Benzo(b)fluoranthene	14.97	252	124357m	6.805	ng	
89) Benzo(k)fluoranthene	15.00	252	40695m	2.331	ng	
90) Benzo(a)pyrene	15.33	252	74390	4.514	ng	97
92) Benzo(a,h,i)perylene	17.23	276	41374	3.243	ng	94

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

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