

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
 Data File : BF111473.D
 Acq On : 15 Dec 2018 17:20
 Operator : JU/SJ
 Sample : J6316-21
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BORING-SAMPLE-STEAL-WELL-YA

Manual Integrations
 APPROVED

Sohil
 12/17/2018 3:12:11 PM

Quant Time: Dec 16 01:36:28 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	194572	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	745775	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	332092	20.00	ng	0.00
64) Phenanthrene-d10	11.31	188	466561	20.00	ng	0.00
76) Chrysene-d12	13.96	240	384851	20.00	ng	0.00
87) Perylene-d12	15.40	264	280996	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	1328521	113.62	ng	0.01
7) Phenol-d6	6.44	99	1703273	109.51	ng	0.00
23) Nitrobenzene-d5	7.36	82	1042519	83.24	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	249888	84.00	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1731646	80.54	ng	0.00
79) Terphenyl-d14	12.90	244	1158359	70.68	ng	0.00
Target Compounds						
10) Phenol	6.45	94	118241	6.825	ng	99
49) Acenaphthylene	9.69	152	101059	3.183	ng	93
50) Dimethylphthalate	9.55	163	184293	7.706	ng	98
66) n-Nitrosodiphenylamine	10.49	169	194257	12.074	ng	97
71) Phenanthrene	11.34	178	524086	21.800	ng	93
72) Anthracene	11.39	178	169140	6.879	ng	95
73) Carbazole	11.54	167	64226	2.526	ng	95
75) Fluoranthene	12.53	202	1090998	46.387	ng	98
78) Pyrene	12.76	202	1055797	38.911	ng	96
81) Benzo(a)anthracene	13.94	228	583079	27.864	ng	97
83) Chrysene	13.98	228	590129m	26.770	ng	
84) Bis(2-ethylhexyl)phthalate	13.94	149	34106	2.019	ng	# 97
86) Indeno(1,2,3-cd)pyrene	16.82	276	245203m	15.412	ng	
88) Benzo(b)fluoranthene	14.99	252	687737m	41.980	ng	
89) Benzo(k)fluoranthene	15.01	252	177909m	11.365	ng	
90) Benzo(a)pyrene	15.34	252	435508	29.480	ng	97
91) Dibenzo(a,h)anthracene	16.82	278	57948m	4.973	ng	
92) Benzo(g,h,i)perylene	17.24	276	248642	21.737	ng	96

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

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