

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121418\
 Data File : BF111431.D
 Acq On : 14 Dec 2018 20:57
 Operator : JU/SJ
 Sample : J6199-05 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-04-121218-A

Manual Integrations
 APPROVED

Sohil
 12/17/2018 2:36:47 PM

Quant Time: Dec 14 22:34:07 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	174609	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	658006	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	260567	20.00	ng	0.00
64) Phenanthrene-d10	11.31	188	366230	20.00	ng	0.00
76) Chrysene-d12	13.95	240	364234	20.00	ng	0.00
87) Perylene-d12	15.39	264	245872	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	490708	46.77	ng	0.00
7) Phenol-d6	6.43	99	612538	43.89	ng	-0.01
23) Nitrobenzene-d5	7.35	82	356020	32.22	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	82311	35.26	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	591108	35.04	ng	0.00
79) Terphenyl-d14	12.90	244	424600	27.37	ng	0.00
Target Compounds						
10) Phenol	6.44	94	39211	2.522	ng	93
50) Dimethylphthalate	9.54	163	46523	2.479	ng	# 98
71) Phenanthrene	11.33	178	41214	2.184	ng	# 90
75) Fluoranthene	12.52	202	75442	4.086	ng	98
78) Pyrene	12.75	202	85243	3.319	ng	98
81) Benzo(a)anthracene	13.94	228	49477	2.498	ng	95
83) Chrysene	13.97	228	44002	2.109	ng	96
88) Benzo(b)fluoranthene	14.97	252	45281m	3.159	ng	
90) Benzo(a)pyrene	15.33	252	32223	2.493	ng	# 86

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

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