

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121018\
 Data File : BF111266.D
 Acq On : 11 Dec 2018 2:16
 Operator : JU/SJ
 Sample : J6197-01 2X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-02-120718

Manual Integrations
 APPROVED

Sohil
 12/11/2018 2:23:49 PM

Quant Time: Dec 11 06:37:43 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 10 01:15:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	443541	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	1575977	20.00	ng	0.00
39) Acenaphthene-d10	9.84	164	601732	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	990642	20.00	ng	0.00
76) Chrysene-d12	13.96	240	692834	20.00	ng	0.00
87) Perylene-d12	15.40	264	535612	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1363803	48.55	ng	0.00
7) Phenol-d6	6.43	99	1684091	48.24	ng	0.00
23) Nitrobenzene-d5	7.36	82	913677	32.52	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	233110	43.73	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	1451211	34.00	ng	0.00
79) Terphenyl-d14	12.91	244	1116655	38.66	ng	0.00
Target Compounds						
10) Phenol	6.44	94	81798	2.004	ng	94
50) Dimethylphthalate	9.55	163	93969	2.116	ng	99
71) Phenanthrene	11.35	178	166626	3.335	ng	93
75) Fluoranthene	12.53	202	220139	4.823	ng	99
78) Pyrene	12.76	202	197496	3.989	ng	96
81) Benzo(a)anthracene	13.95	228	81319	2.104	ng	94
83) Chrysene	13.99	228	82239	2.064	ng	96
88) Benzo(b)fluoranthene	14.99	252	83737m	2.827	ng	
90) Benzo(a)pyrene	15.34	252	56944	2.088	ng	# 74

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

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