

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121418\
 Data File : BF111426.D
 Acq On : 14 Dec 2018 18:40
 Operator : JU/SJ
 Sample : J6319-07DL 5X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 061_SW-N-1DL

Manual Integrations
 APPROVED

Sohil
 12/17/2018 2:36:37 PM

Quant Time: Dec 14 19:09:29 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	171415	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	675982	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	280209	20.00	ng	0.00
64) Phenanthrene-d10	11.31	188	387618	20.00	ng	0.00
76) Chrysene-d12	13.95	240	363148	20.00	ng	0.00
87) Perylene-d12	15.39	264	291080	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.41	112	240524	23.35	ng	0.00
7) Phenol-d6	6.43	99	297931	21.74	ng	-0.01
23) Nitrobenzene-d5	7.35	82	174149	15.34	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	41730	16.62	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	304327	16.78	ng	0.00
79) Terphenyl-d14	12.89	244	192341	12.44	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
52) Acenaphthene	9.86	154	34188	2.019	ng	97
58) Fluorene	10.37	166	45932	2.577	ng	99
71) Phenanthrene	11.33	178	594053	29.742	ng	93
72) Anthracene	11.39	178	152399	7.461	ng	94
75) Fluoranthene	12.52	202	606541	31.041	ng	98
78) Pyrene	12.75	202	750831	29.326	ng	96
81) Benzo(a)anthracene	13.94	228	325475	16.483	ng	99
83) Chrysene	13.97	228	285932	13.746	ng	97
86) Indeno(1,2,3-cd)pyrene	16.80	276	80961	5.393	ng	# 91
88) Benzo(b)fluoranthene	14.98	252	274064m	16.150	ng	
89) Benzo(k)fluoranthene	15.00	252	73691m	4.544	ng	
90) Benzo(a)pyrene	15.33	252	202627	13.241	ng	97
92) Benzo(a,h,i)perylene	17.23	276	86088m	7.265	ng	

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

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