

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121118\
 Data File : BF111275.D
 Acq On : 11 Dec 2018 13:36
 Operator : JU/SJ
 Sample : J6234-01RE 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VNJ-218RE

Manual Integrations
 APPROVED

Sohil
 12/12/2018 5:30:02 PM

Quant Time: Dec 11 17:19:56 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	276029	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	1431127	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	338107	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	471624	20.00	ng	0.00
76) Chrysene-d12	13.95	240	358401	20.00	ng	0.00
87) Perylene-d12	15.39	264	437524	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	536280	30.68	ng	0.00
7) Phenol-d6	6.42	99	657406	30.26	ng	-0.01
23) Nitrobenzene-d5	7.35	82	318110	12.47	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	50080	16.72	ng	-0.01
45) 2-Fluorobiphenyl	9.15	172	592310	22.81	ng	-0.01
79) Terphenyl-d14	12.90	244	206983	13.85	ng	0.00
Target Compounds						
71) Phenanthrene	11.34	178	60227	2.532	ng	95
75) Fluoranthene	12.53	202	97618	4.492	ng	94
78) Pyrene	12.75	202	90066	3.517	ng	96
81) Benzo(a)anthracene	13.94	228	43308	2.166	ng	94
83) Chrysene	13.97	228	44366	2.152	ng	97
88) Benzo(b)fluoranthene	14.97	252	71081m	2.938	ng	
90) Benzo(a)pyrene	15.33	252	47857	2.148	ng	# 80

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

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