

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120518\
 Data File : BF111116.D
 Acq On : 5 Dec 2018 19:26
 Operator : JU/SJ
 Sample : J5764-23 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-113018-B

Manual Integrations
 APPROVED

mohammad
 12/6/2018 12:11:52 PM

Quant Time: Dec 06 01:21:35 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 05 17:30:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	555214	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	2135977	20.00	ng	0.00
39) Acenaphthene-d10	9.84	164	947091	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	1393410	20.00	ng	0.00
76) Chrysene-d12	13.95	240	1052430	20.00	ng	0.00
87) Perylene-d12	15.39	264	772670	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	1565687	44.00	ng	0.00
7) Phenol-d6	6.43	99	2009947	43.85	ng	0.00
23) Nitrobenzene-d5	7.36	82	1186939	32.90	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	300339	39.89	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	2075182	31.51	ng	0.00
79) Terphenyl-d14	12.90	244	1441570	28.05	ng	0.00
Target Compounds						
10) Phenol	6.43	94	113974	2.281	ng	94
50) Dimethylphthalate	9.55	163	156630	2.376	ng	96
71) Phenanthrene	11.34	178	402996	6.013	ng	94
75) Fluoranthene	12.53	202	445489	6.792	ng	97
78) Pyrene	12.76	202	412703	5.033	ng	94
81) Benzo(a)anthracene	13.94	228	189112	3.067	ng	98
83) Chrysene	13.98	228	175305	2.864	ng	95
88) Benzo(b)fluoranthene	14.97	252	164397m	3.755	ng	
90) Benzo(a)pyrene	15.33	252	112753	2.808	ng	# 93

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

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