

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120618\
 Data File : BF111159.D
 Acq On : 7 Dec 2018 00:55
 Operator : JU/SJ
 Sample : J6188-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BU-03-120418-A

Manual Integrations
 APPROVED

mohammad
 12/10/2018 12:19:12 PM

Quant Time: Dec 07 02:49:28 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	518699	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	2043129	20.00	ng	0.00
39) Acenaphthene-d10	9.84	164	954200	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	1386310	20.00	ng	0.00
76) Chrysene-d12	13.96	240	963732	20.00	ng	0.00
87) Perylene-d12	15.39	264	810537	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	2938596	88.40	ng	0.00
7) Phenol-d6	6.43	99	3788247	88.47	ng	0.00
23) Nitrobenzene-d5	7.36	82	2386430	69.15	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	683999	90.18	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	3456638	63.40	ng	0.00
79) Terphenyl-d14	12.91	244	2241680	47.64	ng	0.00
Target Compounds						
10) Phenol	6.45	94	284641	6.097	ng	97
50) Dimethylphthalate	9.56	163	651779	9.815	ng	# 95
71) Phenanthrene	11.35	178	191952	2.879	ng	96
75) Fluoranthene	12.53	202	336848	5.162	ng	93
78) Pyrene	12.76	202	309540	4.123	ng	96
81) Benzo(a)anthracene	13.94	228	183857	3.256	ng	97
83) Chrysene	13.98	228	152708	2.724	ng	94
88) Benzo(b)fluoranthene	14.98	252	208496m	4.540	ng	
90) Benzo(a)pyrene	15.33	252	127800	3.034	ng	# 93

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

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