

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112718\
 Data File : BF111002.D
 Acq On : 27 Nov 2018 11:48
 Operator : JU/SJ
 Sample : J5774-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-02-112118

Manual Integrations
 APPROVED

mohammad
 11/28/2018 4:42:04 PM

Quant Time: Nov 27 13:24:25 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF112618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 27 06:43:50 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	49311	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	186473	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	73334	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	129361	20.00	ng	0.00
76) Chrysene-d12	14.14	240	114326	20.00	ng	0.00
87) Perylene-d12	15.67	264	77515	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	304198	102.30	ng	0.01
7) Phenol-d6	6.57	99	368239	95.03	ng	0.00
23) Nitrobenzene-d5	7.50	82	232053	71.22	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	68692	94.32	ng	0.00
45) 2-Fluorobiphenyl	9.29	172	385473	76.15	ng	0.00
79) Terphenyl-d14	13.06	244	353804	60.24	ng	0.00
Target Compounds						
10) Phenol	6.58	94	20005	4.584	ng	83
50) Dimethylphthalate	9.69	163	53884	10.019	ng	100
71) Phenanthrene	11.50	178	18341	2.677	ng	98
75) Fluoranthene	12.70	202	39166	5.562	ng	97
78) Pyrene	12.93	202	44293	5.305	ng	97
81) Benzo(a)anthracene	14.13	228	22063	3.269	ng	# 94
83) Chrysene	14.16	228	24150m	3.635	ng	
88) Benzo(b)fluoranthene	15.22	252	20575m	4.518	ng	
90) Benzo(a)pyrene	15.62	252	14974	3.533	ng	# 83
92) Benzo(g,h,i)perylene	17.77	276	9532	2.827	ng	# 81

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

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