

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112718\
 Data File : BF110999.D
 Acq On : 27 Nov 2018 10:25
 Operator : JU/SJ
 Sample : J5774-05RE 2X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-03-112118-ARE

Manual Integrations
 APPROVED

mohammad
 11/28/2018 4:41:53 PM

Quant Time: Nov 27 13:00:23 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	54654	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	204102	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	75419	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	121838	20.00	ng	0.00
76) Chrysene-d12	14.13	240	131552	20.00	ng	0.00
87) Perylene-d12	15.67	264	115395	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	163076	49.48	ng	0.00
7) Phenol-d6	6.56	99	206555	48.09	ng	0.00
23) Nitrobenzene-d5	7.50	82	127996	35.89	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	31473	42.02	ng	0.00
45) 2-Fluorobiphenyl	9.29	172	200724	38.56	ng	0.00
79) Terphenyl-d14	13.06	244	165636	24.51	ng	0.00
Target Compounds						
10) Phenol	6.58	94	11209	2.317	ng	92
50) Dimethylphthalate	9.69	163	25204	4.557	ng	99
52) Acenaphthene	10.01	154	12614	2.745	ng	96
58) Fluorene	10.53	166	21445	4.111	ng	96
71) Phenanthrene	11.50	178	120743	18.711	ng	99
72) Anthracene	11.56	178	35164	5.351	ng	97
75) Fluoranthene	12.70	202	151951	22.911	ng	99
78) Pyrene	12.93	202	165035	17.176	ng	97
81) Benzo(a)anthracene	14.12	228	79392	10.222	ng	98
83) Chrysene	14.16	228	73405m	9.601	ng	
86) Indeno(1,2,3-cd)pyrene	17.27	276	23378	4.260	ng	94
88) Benzo(b)fluoranthene	15.22	252	67517m	9.960	ng	
89) Benzo(k)fluoranthene	15.24	252	25635m	3.674	ng	
90) Benzo(a)pyrene	15.61	252	52839	8.374	ng	# 96
92) Benzo(a,h,i)perylene	17.76	276	23201	4.622	ng	96

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

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