

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102518\
 Data File : BF110182.D
 Acq On : 26 Oct 2018 1:13
 Operator : JU/SJ
 Sample : J5200-15 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-02-102318-B

Manual Integrations
 APPROVED

Sohil
 10/26/2018 3:49:31 PM

Quant Time: Oct 26 07:14:16 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 23 15:06:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	138454	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	517193	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	203007	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	347575	20.00	ng	-0.01
76) Chrysene-d12	14.15	240	245636	20.00	ng	-0.01
87) Perylene-d12	15.68	264	186486	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	387125	45.53	ng	-0.01
7) Phenol-d6	6.59	99	476202	43.79	ng	-0.02
23) Nitrobenzene-d5	7.53	82	273858	31.41	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	79364	39.47	ng	-0.02
45) 2-Fluorobiphenyl	9.32	172	484103	37.45	ng	-0.02
79) Terphenyl-d14	13.09	244	392317	33.22	ng	-0.02
Target Compounds						
10) Phenol	6.60	94	27721	2.385	ng	Qvalue # 62
50) Dimethylphthalate	9.72	163	61254	4.088	ng	99
71) Phenanthrene	11.53	178	186617	10.261	ng	98
72) Anthracene	11.58	178	52106	2.858	ng	99
75) Fluoranthene	12.72	202	280348	15.189	ng	97
78) Pyrene	12.95	202	255615	14.515	ng	99
81) Benzo(a)anthracene	14.14	228	109618	7.525	ng	95
83) Chrysene	14.17	228	96416	6.969	ng	96
86) Indeno(1,2,3-cd)pyrene	17.24	276	42594	3.711	ng	99
88) Benzo(b)fluoranthene	15.22	252	113266m	10.518	ng	
89) Benzo(k)fluoranthene	15.25	252	28769m	2.717	ng	
90) Benzo(a)pyrene	15.61	252	73015	7.368	ng	# 96
92) Benzo(a,h,i)perylene	17.72	276	44497	5.281	ng	# 86

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

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