

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111918\
 Data File : BF110861.D
 Acq On : 19 Nov 2018 18:53
 Operator : JU/SJ
 Sample : J5774-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-04-111618

Manual Integrations
 APPROVED

Sohil
 11/20/2018 3:48:56 PM

Quant Time: Nov 20 10:54:59 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	56208	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	214390	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	88149	20.00	ng	-0.01
64) Phenanthrene-d10	11.48	188	141716	20.00	ng	0.00
76) Chrysene-d12	14.14	240	112250	20.00	ng	0.00
87) Perylene-d12	15.67	264	73993	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	348104	100.60	ng	0.00
7) Phenol-d6	6.57	99	432993	97.85	ng	0.00
23) Nitrobenzene-d5	7.50	82	264357	71.01	ng	-0.01
42) 2,4,6-Tribromophenol	10.78	330	76935	86.62	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	439291	71.17	ng	0.00
79) Terphenyl-d14	13.07	244	381422	63.64	ng	0.00
Target Compounds						
10) Phenol	6.58	94	19168	3.736	ng	Qvalue 81
50) Dimethylphthalate	9.69	163	37928	5.766	ng	# 97
71) Phenanthrene	11.50	178	83277	10.968	ng	99
72) Anthracene	11.56	178	25954m	3.389	ng	
75) Fluoranthene	12.70	202	157402	20.678	ng	99
78) Pyrene	12.94	202	168442	19.489	ng	97
81) Benzo(a)anthracene	14.13	228	76312	11.374	ng	96
83) Chrysene	14.16	228	83098m	13.000	ng	
86) Indeno(1,2,3-cd)pyrene	17.26	276	23181	5.054	ng	98
88) Benzo(b)fluoranthene	15.22	252	63384m	14.319	ng	
89) Benzo(k)fluoranthene	15.24	252	20615m	4.713	ng	
90) Benzo(a)pyrene	15.61	252	43494	10.815	ng	# 87
91) Dibenzo(a,h)anthracene	17.26	278	8059	2.368	ng	# 41
92) Benzo(g,h,i)perylene	17.74	276	26359	7.806	ng	# 91

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

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