

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071118\
 Data File : BF107297.D
 Acq On : 11 Jul 2018 13:59
 Operator : JU/SJ
 Sample : J3933-12
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP-12-W1

Manual Integrations
 APPROVED

Sohil
 7/12/2018 2:19:11 PM

Quant Time: Jul 11 15:40:24 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 29 14:38:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	80278	20.00	ng	-0.03
21) Naphthalene-d8	8.21	136	302493	20.00	ng	-0.04
38) Acenaphthene-d10	9.98	164	148696	20.00	ng	-0.03
63) Phenanthrene-d10	11.48	188	263334	20.00	ng	-0.02
75) Chrysene-d12	14.14	240	223572	20.00	ng	-0.02
86) Perylene-d12	15.69	264	254787	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	447517	94.40	ng	-0.03
7) Phenol-d6	6.58	99	561133	94.78	ng	-0.03
23) Nitrobenzene-d5	7.50	82	368565	67.71	ng	-0.04
41) 2,4,6-Tribromophenol	10.78	330	167719	97.32	ng	-0.02
44) 2-Fluorobiphenyl	9.29	172	633216	69.37	ng	-0.03
78) Terphenyl-d14	13.07	244	589241	63.84	ng	-0.02
Target Compounds						
10) Phenol	6.60	94	29905	4.426	ng	99
20) 3+4-Methylphenols	7.38	107	481	Below Cal	#	70
48) Acenaphthylene	9.84	152	66701	4.529	ng	99
49) Dimethylphthalate	9.68	163	179391	16.085	ng	98
70) Phenanthrene	11.50	178	218186	17.178	ng	99
71) Anthracene	11.55	178	54726	4.221	ng	98
72) Carbazole	11.71	167	25021	2.061	ng	96
74) Fluoranthene	12.70	202	585974	41.858	ng	96
77) Pyrene	12.94	202	557317	37.802	ng	100
80) Benzo(a)anthracene	14.13	228	381196	27.975	ng	95
82) Chrysene	14.17	228	325412	25.958	ng	98
83) Bis(2-ethylhexyl)phthalate	14.08	149	313999	40.518	ng	# 97
85) Indeno(1,2,3-cd)pyrene	17.28	276	236783	22.258	ng	# 85
87) Benzo(b)fluoranthene	15.23	252	565907m	35.755	ng	
88) Benzo(k)fluoranthene	15.25	252	161358m	10.706	ng	
89) Benzo(a)pyrene	15.62	252	388476m	27.610	ng	
90) Dibenzo(a,h)anthracene	17.28	278	62404m	5.233	ng	
91) Benzo(g,h,i)perylene	17.77	276	213774	18.342	ng	# 89

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

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