

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010919\
 Data File : BF111974.D
 Acq On : 10 Jan 2019 3:27
 Operator : JU/SJ
 Sample : K1044-05DL 10X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CPSB-10(10-15)DL

Manual Integrations
 APPROVED

Sohil
 1/10/2019 11:46:15 AM

Quant Time: Jan 10 07:16:26 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 07 14:33:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	116530	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	446203	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	162417	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	282848	20.00	ng	0.00
76) Chrysene-d12	14.06	240	193151	20.00	ng	0.00
87) Perylene-d12	15.55	264	152596	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	58247	8.55	ng	0.00
7) Phenol-d6	6.52	99	75560	8.61	ng	-0.01
23) Nitrobenzene-d5	7.44	82	41976	5.03	ng	-0.02
42) 2,4,6-Tribromophenol	10.72	330	15091	7.14	ng	0.00
45) 2-Fluorobiphenyl	9.24	172	72389	6.50	ng	-0.01
79) Terphenyl-d14	13.00	244	64499	6.34	ng	0.00
Target Compounds						
52) Acenaphthene	9.96	154	34618	3.406	ng	Qvalue 95
58) Fluorene	10.47	166	51420	4.781	ng	94
71) Phenanthrene	11.44	178	354612	23.449	ng	99
72) Anthracene	11.49	178	104649	6.817	ng	99
75) Fluoranthene	12.63	202	336144	21.652	ng	98
78) Pyrene	12.86	202	433453	32.615	ng	99
81) Benzo(a)anthracene	14.05	228	156370	14.067	ng	99
83) Chrysene	14.09	228	141952	13.393	ng	98
86) Indeno(1,2,3-cd)pyrene	17.05	276	44573	5.768	ng	97
88) Benzo(b)fluoranthene	15.12	252	111418m	12.206	ng	
89) Benzo(k)fluoranthene	15.14	252	30724m	3.581	ng	
90) Benzo(a)pyrene	15.49	252	85858	10.697	ng	# 96
91) Dibenzo(a,h)anthracene	17.06	278	15356	2.379	ng	# 80
92) Benzo(g,h,i)perylene	17.51	276	45082	7.140	ng	96

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

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