

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102418\
 Data File : BF110125.D
 Acq On : 24 Oct 2018 19:55
 Operator : JU/SJ
 Sample : J5198-29 2X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-102318-E

Manual Integrations
 APPROVED

Sohil
 10/25/2018 11:15:27 AM

Quant Time: Oct 25 04:31:41 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	136225	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	506118	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	197969	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	343349	20.00	ng	-0.01
76) Chrysene-d12	14.15	240	259691	20.00	ng	-0.01
87) Perylene-d12	15.68	264	185480	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	416800	49.82	ng	-0.02
7) Phenol-d6	6.59	99	501243	46.85	ng	-0.02
23) Nitrobenzene-d5	7.53	82	273104	32.01	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	83405	42.53	ng	-0.01
45) 2-Fluorobiphenyl	9.32	172	502495	39.86	ng	-0.02
79) Terphenyl-d14	13.09	244	448933	35.95	ng	-0.02
Target Compounds						
10) Phenol	6.60	94	65959	5.767	ng	Qvalue # 60
50) Dimethylphthalate	9.72	163	68791	4.707	ng	98
71) Phenanthrene	11.53	178	131396	7.314	ng	100
72) Anthracene	11.58	178	42446	2.357	ng	97
75) Fluoranthene	12.72	202	276336	15.156	ng	99
78) Pyrene	12.95	202	261900	14.067	ng	98
81) Benzo(a)anthracene	14.14	228	118379	7.687	ng	95
83) Chrysene	14.18	228	110534	7.557	ng	98
84) Bis(2-ethylhexyl)phthalate	14.11	149	48941	4.480	ng	96
86) Indeno(1,2,3-cd)pyrene	17.24	276	45158	3.721	ng	# 91
88) Benzo(b)fluoranthene	15.23	252	120082m	11.211	ng	
89) Benzo(k)fluoranthene	15.25	252	38237m	3.631	ng	
90) Benzo(a)pyrene	15.62	252	76155	7.727	ng	# 93
92) Benzo(a,h,i)perylene	17.72	276	44759	5.341	ng	# 85

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

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