

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083118\
 Data File : BF108702.D
 Acq On : 31 Aug 2018 21:11
 Operator : JU/SJ
 Sample : J4720-04 4X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BZTRANS-4

Manual Integrations
 APPROVED

Sohil
 9/4/2018 3:53:33 PM

Quant Time: Sep 04 15:39:10 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 30 15:06:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.01	152	59933	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	253383	20.00	ng	0.00
38) Acenaphthene-d10	10.04	164	108598	20.00	ng	0.00
63) Phenanthrene-d10	11.55	188	196406	20.00	ng	0.00
75) Chrysene-d12	14.21	240	204877	20.00	ng	0.00
86) Perylene-d12	15.76	264	144677	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.78	112	43015m	12.47	ng	0.13
7) Phenol-d6	6.67	99	86644m	17.53	ng	0.02
23) Nitrobenzene-d5	7.59	82	56076	11.23	ng	0.02
41) 2,4,6-Tribromophenol	10.85	330	20500	14.30	ng	0.00
44) 2-Fluorobiphenyl	9.35	172	114197	14.14	ng	0.00
78) Terphenyl-d14	13.14	244	121797	11.55	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.74	163	37369	4.370	ng	# 94
70) Phenanthrene	11.57	178	83463	8.419	ng	# 74
74) Fluoranthene	12.78	202	111914	10.029	ng	# 79
77) Pyrene	13.01	202	86808	5.636	ng	# 81
82) Chrysene	14.24	228	39524	3.291	ng	# 90
83) Bis(2-ethylhexyl)phthalate	14.14	149	21178	2.544	ng	# 91
85) Indeno(1,2,3-cd)pyrene	17.38	276	18594	2.226	ng	# 91
87) Benzo(b)fluoranthene	15.29	252	38757m	4.361	ng	# 84
89) Benzo(a)pyrene	15.69	252	22318	2.779	ng	# 84
91) Benzo(g,h,i)perylene	17.88	276	18286	2.933	ng	# 87

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

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