

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083118\
 Data File : BF108694.D
 Acq On : 31 Aug 2018 17:35
 Operator : JU/SJ
 Sample : J4720-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BZTRANS-3

Manual Integrations
 APPROVED

Sohil
 9/4/2018 3:53:14 PM

Quant Time: Sep 04 15:31:55 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	7.00	152	69232	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	277521	20.00	ng	0.00
38) Acenaphthene-d10	10.04	164	119879	20.00	ng	0.00
63) Phenanthrene-d10	11.54	188	231582	20.00	ng	0.00
75) Chrysene-d12	14.20	240	221769	20.00	ng	0.00
86) Perylene-d12	15.75	264	162725	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	187615	47.08	ng	0.00
7) Phenol-d6	6.64	99	366693	64.24	ng	0.00
23) Nitrobenzene-d5	7.57	82	235564	43.08	ng	0.00
41) 2,4,6-Tribromophenol	10.84	330	92906	58.71	ng	-0.01
44) 2-Fluorobiphenyl	9.35	172	477102	53.51	ng	0.00
78) Terphenyl-d14	13.12	244	470977	41.26	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.74	163	163551	17.325	ng	99
70) Phenanthrene	11.57	178	157607m	13.483	ng	
74) Fluoranthene	12.76	202	236850	18.001	ng	94
77) Pyrene	12.99	202	188029	11.279	ng	98
80) Benzo(a)anthracene	14.18	228	58452	4.324	ng	97
82) Chrysene	14.22	228	90533	6.963	ng	98
83) Bis(2-ethylhexyl)phthalate	14.13	149	32016	3.553	ng	100
85) Indeno(1,2,3-cd)pyrene	17.38	276	32005	3.540	ng	# 85
87) Benzo(b)fluoranthene	15.29	252	91692m	9.173	ng	
88) Benzo(k)fluoranthene	15.31	252	36636m	3.700	ng	
89) Benzo(a)pyrene	15.69	252	55866	6.184	ng	# 90
91) Benzo(a,h,i)perylene	17.87	276	33973	4.845	ng	91

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

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