

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083118\
 Data File : BF108697.D
 Acq On : 31 Aug 2018 18:56
 Operator : JU/SJ
 Sample : J4720-02 5X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BZTRANS-2

Manual Integrations
 APPROVED

Sohil
 9/4/2018 3:53:21 PM

Quant Time: Sep 01 04:05:24 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	60132	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	285427	20.00	ng	0.00
38) Acenaphthene-d10	10.04	164	119292	20.00	ng	0.00
63) Phenanthrene-d10	11.54	188	214272	20.00	ng	0.00
75) Chrysene-d12	14.19	240	222173	20.00	ng	0.00
86) Perylene-d12	15.75	264	165429	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.90	112	20691m	5.98	ng	0.25
7) Phenol-d6	6.71	99	54905m	11.07	ng	0.06
23) Nitrobenzene-d5	7.60	82	44556m	7.92	ng	0.03
41) 2,4,6-Tribromophenol	10.85	330	13628	8.65	ng	0.00
44) 2-Fluorobiphenyl	9.35	172	88387	9.96	ng	0.00
78) Terphenyl-d14	13.12	244	88716	7.76	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.75	163	29459	3.136	ng	Qvalue # 99
70) Phenanthrene	11.57	178	210568	19.469	ng	99
74) Fluoranthene	12.77	202	258537	21.237	ng	94
77) Pyrene	12.99	202	197267	11.812	ng	98
80) Benzo(a)anthracene	14.19	228	44432	3.281	ng	96
82) Chrysene	14.22	228	97880	7.515	ng	98
85) Indeno(1,2,3-cd)pyrene	17.38	276	35214	3.888	ng	# 87
87) Benzo(b)fluoranthene	15.28	252	79428m	7.816	ng	
88) Benzo(k)fluoranthene	15.30	252	50008m	4.968	ng	
89) Benzo(a)pyrene	15.69	252	56170	6.116	ng	96
91) Benzo(g,h,i)perylene	17.87	276	29957	4.202	ng	# 91

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

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