

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081618\
 Data File : BF108207.D
 Acq On : 16 Aug 2018 20:50
 Operator : JU/SJ
 Sample : J4499-07 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LSB-07(0-1)

Manual Integrations
 APPROVED

Sohil
 8/17/2018 3:30:52 PM

Quant Time: Aug 17 04:36:02 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 08 12:24:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.02	152	156320	20.00	ng	0.00
21) Naphthalene-d8	8.30	136	570454	20.00	ng	0.00
38) Acenaphthene-d10	10.06	164	248639	20.00	ng	0.00
63) Phenanthrene-d10	11.55	188	427803	20.00	ng	0.00
75) Chrysene-d12	14.21	240	420634	20.00	ng	0.00
86) Perylene-d12	15.77	264	316706	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	347915	40.29	ng	0.00
7) Phenol-d6	6.62	99	440718	38.41	ng	0.00
23) Nitrobenzene-d5	7.57	82	285818	27.96	ng	0.00
41) 2,4,6-Tribromophenol	10.85	330	88236	35.58	ng	0.00
44) 2-Fluorobiphenyl	9.37	172	474564	30.64	ng	0.00
78) Terphenyl-d14	13.14	244	433819	26.22	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.76	163	41792	2.413	ng	Qvalue # 97
70) Phenanthrene	11.58	178	244071	12.096	ng	97
71) Anthracene	11.63	178	47898	2.316	ng	99
74) Fluoranthene	12.78	202	746804	33.912	ng	95
77) Pyrene	13.01	202	618235	23.215	ng	99
80) Benzo(a)anthracene	14.20	228	349714	14.487	ng	99
82) Chrysene	14.24	228	316969	14.322	ng	99
85) Indeno(1,2,3-cd)pyrene	17.38	276	120262	6.271	ng	# 89
87) Benzo(b)fluoranthene	15.30	252	366404m	19.361	ng	
88) Benzo(k)fluoranthene	15.33	252	150034m	8.625	ng	
89) Benzo(a)pyrene	15.70	252	226195	13.348	ng	96
90) Dibenzo(a,h)anthracene	17.39	278	31945m	2.278	ng	
91) Benzo(a,h,i)perylene	17.88	276	114099	8.264	ng	# 91

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

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