

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081518\
 Data File : BF108178.D
 Acq On : 15 Aug 2018 21:06
 Operator : JU/SJ
 Sample : J4460-11
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RA-081018-FL-011

Manual Integrations
 APPROVED

Sohil
 8/16/2018 5:27:06 PM

Quant Time: Aug 16 03:48:52 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	216805	20.00	ng	0.00
21) Naphthalene-d8	8.30	136	819688	20.00	ng	0.00
38) Acenaphthene-d10	10.06	164	386556	20.00	ng	0.00
63) Phenanthrene-d10	11.56	188	621014	20.00	ng	0.00
75) Chrysene-d12	14.21	240	445159	20.00	ng	0.00
86) Perylene-d12	15.77	264	459987	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.65	112	1262499	105.43	ng	0.02
7) Phenol-d6	6.64	99	1696065	106.58	ng	0.01
23) Nitrobenzene-d5	7.58	82	1161481	79.07	ng	0.00
41) 2,4,6-Tribromophenol	10.85	330	356033	92.34	ng	0.00
44) 2-Fluorobiphenyl	9.38	172	1967260	81.71	ng	0.00
78) Terphenyl-d14	13.15	244	1463533	83.59	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.76	163	202239	7.511	ng	Qvalue # 99
70) Phenanthrene	11.58	178	267814	9.143	ng	98
71) Anthracene	11.63	178	75082	2.501	ng	98
74) Fluoranthene	12.78	202	302651	9.467	ng	94
77) Pyrene	13.01	202	260276	9.235	ng	98
80) Benzo(a)anthracene	14.20	228	148638	5.818	ng	98
82) Chrysene	14.24	228	117978	5.037	ng	96
85) Indeno(1,2,3-cd)pyrene	17.38	276	52769	2.600	ng	# 86
87) Benzo(b)fluoranthene	15.30	252	152088m	5.533	ng	
88) Benzo(k)fluoranthene	15.32	252	55978m	2.216	ng	
89) Benzo(a)pyrene	15.70	252	113818	4.624	ng	98
91) Benzo(a,h,i)perylene	17.88	276	44451	2.217	ng	92

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

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