

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082218\
 Data File : BF108392.D
 Acq On : 22 Aug 2018 19:22
 Operator : JU/SJ
 Sample : J4559-09
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RA-081718-WSW-20-053

Manual Integrations
 APPROVED

Sohil
 8/23/2018 3:43:18 PM

Quant Time: Aug 23 05:05:53 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 22 11:01:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.00	152	145114	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	547510	20.00	ng	0.00
38) Acenaphthene-d10	10.05	164	260087	20.00	ng	0.00
63) Phenanthrene-d10	11.54	188	434982	20.00	ng	0.00
75) Chrysene-d12	14.20	240	251321	20.00	ng	0.00
86) Perylene-d12	15.77	264	284347	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	787996	98.31	ng	0.00
7) Phenol-d6	6.62	99	1082803	101.66	ng	0.00
23) Nitrobenzene-d5	7.57	82	711952	72.56	ng	0.00
41) 2,4,6-Tribromophenol	10.84	330	275019	106.01	ng	0.00
44) 2-Fluorobiphenyl	9.36	172	1076017	66.42	ng	0.00
78) Terphenyl-d14	13.13	244	687811	69.58	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.75	163	370613	20.456	ng	Qvalue # 99
51) Acenaphthene	10.08	154	44166	3.009	ng	96
57) Fluorene	10.60	166	43845	2.605	ng	97
70) Phenanthrene	11.57	178	460528	22.447	ng	100
71) Anthracene	11.62	178	109518	5.208	ng	97
72) Carbazole	11.77	167	51804	2.773	ng	# 92
74) Fluoranthene	12.77	202	649264	28.996	ng	96
77) Pyrene	13.00	202	568964	35.759	ng	99
80) Benzo(a)anthracene	14.19	228	269055	18.654	ng	96
82) Chrysene	14.22	228	248444	18.789	ng	97
83) Bis(2-ethylhexyl)phthalate	14.15	149	24610	2.876	ng	# 79
85) Indeno(1,2,3-cd)pyrene	17.38	276	191888	16.748	ng	# 86
87) Benzo(b)fluoranthene	15.29	252	374207m	22.024	ng	
88) Benzo(k)fluoranthene	15.32	252	142579m	9.129	ng	
89) Benzo(a)pyrene	15.70	252	275705	18.121	ng	# 95
90) Dibenzo(a,h)anthracene	17.38	278	45802	3.638	ng	# 62
91) Benzo(a,h,i)perylene	17.88	276	182909	14.755	ng	# 89

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

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