

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082418\
 Data File : BF108480.D
 Acq On : 24 Aug 2018 22:57
 Operator : JU/SJ
 Sample : J4581-18
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP1-Q3-D2

Manual Integrations
 APPROVED

Sohil
 8/27/2018 12:56:52 PM

Quant Time: Aug 25 01:43:41 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	125675	20.00	ng	-0.01
21) Naphthalene-d8	8.28	136	481178	20.00	ng	0.00
38) Acenaphthene-d10	10.04	164	235651	20.00	ng	-0.01
63) Phenanthrene-d10	11.54	188	429635	20.00	ng	-0.01
75) Chrysene-d12	14.19	240	265082	20.00	ng	-0.01
86) Perylene-d12	15.74	264	277191	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	560777	80.78	ng	0.00
7) Phenol-d6	6.62	99	765511	82.99	ng	-0.01
23) Nitrobenzene-d5	7.56	82	506460	58.73	ng	-0.01
41) 2,4,6-Tribromophenol	10.84	330	206863	88.01	ng	-0.01
44) 2-Fluorobiphenyl	9.35	172	812454	55.35	ng	-0.01
78) Terphenyl-d14	13.12	244	680824	65.30	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.75	163	364677	22.216	ng	99
57) Fluorene	10.59	166	41135	2.697	ng	99
70) Phenanthrene	11.56	178	347600	17.153	ng	98
71) Anthracene	11.61	178	86740	4.176	ng	98
74) Fluoranthene	12.75	202	354186	16.015	ng	96
77) Pyrene	12.99	202	375043	22.347	ng	100
80) Benzo(a)anthracene	14.18	228	146051	9.600	ng	98
82) Chrysene	14.21	228	134238	9.625	ng	97
85) Indeno(1,2,3-cd)pyrene	17.35	276	92904	7.688	ng	# 89
87) Benzo(b)fluoranthene	15.28	252	192761m	11.638	ng	
88) Benzo(k)fluoranthene	15.30	252	56446m	3.707	ng	
89) Benzo(a)pyrene	15.68	252	156871	10.577	ng	98
90) Dibenzo(a,h)anthracene	17.35	278	24640	2.008	ng	95
91) Benzo(g,h,i)perylene	17.84	276	92202m	7.630	ng	

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

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