

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082418\
 Data File : BF108468.D
 Acq On : 24 Aug 2018 17:32
 Operator : JU/SJ
 Sample : J4581-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP1-Q3-A3

Manual Integrations
 APPROVED

Sohil
 8/27/2018 12:56:40 PM

Quant Time: Aug 24 18:41:13 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	132347	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	509609	20.00	ng	0.00
38) Acenaphthene-d10	10.04	164	259139	20.00	ng	-0.01
63) Phenanthrene-d10	11.54	188	503959	20.00	ng	0.00
75) Chrysene-d12	14.19	240	299170	20.00	ng	0.00
86) Perylene-d12	15.75	264	286259	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	745950	102.04	ng	0.00
7) Phenol-d6	6.62	99	1023957	105.41	ng	0.00
23) Nitrobenzene-d5	7.56	82	678336	74.28	ng	-0.01
41) 2,4,6-Tribromophenol	10.84	330	315725	122.14	ng	0.00
44) 2-Fluorobiphenyl	9.36	172	1210560	75.00	ng	0.00
78) Terphenyl-d14	13.12	244	1198157	101.83	ng	0.00
Target Compounds						
10) Phenol	6.64	94	21186	2.108	ng	90
49) Dimethylphthalate	9.75	163	344107	19.063	ng	99
70) Phenanthrene	11.57	178	145654	6.128	ng	97
74) Fluoranthene	12.76	202	202590	7.809	ng	94
77) Pyrene	12.99	202	179208	9.462	ng	99
80) Benzo(a)anthracene	14.18	228	69404	4.042	ng	98
82) Chrysene	14.22	228	68213	4.334	ng	95
85) Indeno(1,2,3-cd)pyrene	17.35	276	38377	2.814	ng	# 90
87) Benzo(b)fluoranthene	15.28	252	77421m	4.526	ng	
89) Benzo(a)pyrene	15.68	252	56248	3.672	ng	97
91) Benzo(g,h,i)perylene	17.85	276	37245	2.985	ng	94

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

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