

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF020119\
 Data File : BF112358.D
 Acq On : 1 Feb 2019 18:18
 Operator : JU/SJ
 Sample : K1297-05
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 A0C17-SS2

Manual Integrations
 APPROVED

apatel
 2/4/2019 12:43:22 PM

Quant Time: Feb 02 00:31:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 28 10:38:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	166029	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	601164	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	262908	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	412025	20.00	ng	-0.01
76) Chrysene-d12	14.00	240	360325	20.00	ng	-0.01
87) Perylene-d12	15.47	264	375392	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	932395	119.28	ng	0.00
7) Phenol-d6	6.49	99	1127333	103.79	ng	0.00
23) Nitrobenzene-d5	7.40	82	725352	69.52	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	276911	90.49	ng	-0.02
45) 2-Fluorobiphenyl	9.19	172	1164895	62.67	ng	-0.01
79) Terphenyl-d14	12.94	244	812290	42.46	ng	-0.01
Target Compounds						
10) Phenol	6.50	94	25377	2.130	ng	83
50) Dimethylphthalate	9.58	163	96673	4.733	ng	98
71) Phenanthrene	11.38	178	84077	3.925	ng	98
75) Fluoranthene	12.57	202	132495	5.767	ng	96
78) Pyrene	12.80	202	112800	4.522	ng	99
81) Benzo(a)anthracene	13.99	228	67327	3.179	ng	98
83) Chrysene	14.03	228	67426	3.335	ng	96
88) Benzo(b)fluoranthene	15.04	252	91798m	4.089	ng	
90) Benzo(a)pyrene	15.40	252	62113	3.075	ng	# 94
92) Benzo(g,h,i)perylene	17.39	276	32640	2.125	ng	94

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

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