

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF013019\
 Data File : BF112327.D
 Acq On : 30 Jan 2019 17:46
 Operator : JU/SJ
 Sample : K1282-25
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-7

Manual Integrations
 APPROVED

mohammad
 2/1/2019 12:06:30 AM

Quant Time: Jan 31 00:33:43 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	169020	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	635463	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	293879	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	503299	20.00	ng	-0.01
76) Chrysene-d12	14.00	240	356968	20.00	ng	-0.01
87) Perylene-d12	15.47	264	359350	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	880962	110.71	ng	0.00
7) Phenol-d6	6.49	99	1089362	98.52	ng	0.00
23) Nitrobenzene-d5	7.40	82	705883	64.01	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	266902	78.03	ng	-0.02
45) 2-Fluorobiphenyl	9.19	172	1173123	56.46	ng	-0.01
79) Terphenyl-d14	12.94	244	851597	44.93	ng	-0.01
Target Compounds						
10) Phenol	6.50	94	57240	4.718	ng	78
50) Dimethylphthalate	9.59	163	96646	4.233	ng	99
71) Phenanthrene	11.38	178	121622	4.648	ng	96
75) Fluoranthene	12.57	202	165669	5.903	ng	98
78) Pyrene	12.80	202	152115	6.155	ng	98
81) Benzo(a)anthracene	13.99	228	73815	3.518	ng	98
83) Chrysene	14.03	228	69957	3.492	ng	97
86) Indeno(1,2,3-cd)pyrene	16.94	276	33365	2.102	ng	# 92
88) Benzo(b)fluoranthene	15.04	252	92252m	4.293	ng	
90) Benzo(a)pyrene	15.40	252	62757	3.245	ng	# 97
92) Benzo(g,h,i)perylene	17.39	276	34832	2.369	ng	99

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

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