

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090419\
 Data File : BF116820.D
 Acq On : 4 Sep 2019 17:21
 Operator : HP/JU
 Sample : K4639-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DW-02

Quant Time: Sep 05 04:06:01 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 30 20:02:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	59899	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	240792	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	122334	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	185894	20.00	ng	0.00
76) Chrysene-d12	14.00	240	143916	20.00	ng	0.00
87) Perylene-d12	15.46	264	127439	20.00	ng	0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	369632	99.97	ng	0.00
7) Phenol-d6	6.48	99	498905	102.76	ng	0.00
23) Nitrobenzene-d5	7.40	82	304343	72.15	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	97117	118.75	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	540557	74.54	ng	0.00
79) Terphenyl-d14	12.95	244	461099	59.27	ng	0.00

Target Compounds

						Qvalue
10) Phenol	6.49	94	15657	2.912	ng	97
50) Dimethylphthalate	9.59	163	82017	9.711	ng	99
71) Phenanthrene	11.39	178	43111	4.166	ng	97
75) Fluoranthene	12.58	202	91325	8.980	ng	99
78) Pyrene	12.80	202	78428	6.130	ng	97
81) Benzo(a)anthracene	13.99	228	20418m	2.242	ng	
83) Chrysene	14.03	228	50169	5.346	ng	# 95
84) Bis(2-ethylhexyl)phthalate	13.99	149	262225	34.145	ng	97
86) Indeno(1,2,3-cd)pyrene	16.92	276	15697	2.083	ng	96
88) Benzo(b)fluoranthene	15.04	252	49563m	6.433	ng	
90) Benzo(a)pyrene	15.40	252	18758	2.798	ng	# 1
92) Benzo(a,h,i)perylene	17.36	276	17277	2.887	ng	# 74

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

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