

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\
 Data File : BF117582.D
 Acq On : 29 Oct 2019 11:24
 Operator : JU
 Sample : K5503-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 176-28-(0-2.5)

Manual Integrations
 APPROVED

mohammad
 10/30/2019 4:52:38 PM

Quant Time: Oct 30 02:28:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 18:53:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	48780	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	198772	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	90754	20.00	ng	-0.01
64) Phenanthrene-d10	11.25	188	113804	20.00	ng	0.00
76) Chrysene-d12	13.89	240	85282	20.00	ng	0.00
87) Perylene-d12	15.33	264	98239	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	346871	105.79	ng	0.00
7) Phenol-d6	6.39	99	475505	107.97	ng	0.00
23) Nitrobenzene-d5	7.29	82	295514	73.40	ng	-0.01
42) 2,4,6-Tribromophenol	10.56	330	68076	86.69	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	448197	69.58	ng	-0.01
79) Terphenyl-d14	12.83	244	255851	48.93	ng	0.00
Target Compounds						
10) Phenol	6.40	94	16514	3.674	ng	Qvalue 83
50) Dimethylphthalate	9.48	163	47717	7.375	ng	99
71) Phenanthrene	11.27	178	33581	5.082	ng	97
75) Fluoranthene	12.46	202	58444	8.835	ng	99
78) Pyrene	12.69	202	57168	7.804	ng	98
81) Benzo(a)anthracene	13.88	228	30485	5.299	ng	96
83) Chrysene	13.92	228	30532	5.415	ng	96
86) Indeno(1,2,3-cd)pyrene	16.74	276	18454	4.384	ng	# 92
88) Benzo(b)fluoranthene	14.92	252	48098m	8.291	ng	
89) Benzo(k)fluoranthene	14.94	252	12140m	2.042	ng	
90) Benzo(a)pyrene	15.27	252	32318	6.127	ng	# 91
92) Benzo(a,h,i)perylene	17.17	276	17565m	4.071	ng	

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

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