

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080219\
 Data File : BF115953.D
 Acq On : 2 Aug 2019 18:18
 Operator : HP/JU
 Sample : K4131-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 1S-D1-D4-COMP-(30)

Manual Integrations
 APPROVED

mohammad
 8/8/2019 4:44:41 PM

Quant Time: Aug 03 00:35:51 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 31 15:31:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	101537	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	400244	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	220488	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	402240	20.00	ng	0.00
76) Chrysene-d12	14.03	240	308777	20.00	ng	0.00
87) Perylene-d12	15.50	264	349668	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	581751	95.91	ng	0.00
7) Phenol-d6	6.48	99	751801	98.24	ng	0.00
23) Nitrobenzene-d5	7.41	82	482294	69.56	ng	0.00
42) 2,4,6-Tribromophenol	10.68	330	199350	93.25	ng	0.00
45) 2-Fluorobiphenyl	9.21	172	842054	71.53	ng	0.00
79) Terphenyl-d14	12.97	244	914605	62.41	ng	0.00
Target Compounds						
10) Phenol	6.49	94	40270	4.469	ng	94
50) Dimethylphthalate	9.60	163	72730	4.845	ng	99
71) Phenanthrene	11.40	178	273535	13.302	ng	99
72) Anthracene	11.46	178	78293	3.762	ng	97
75) Fluoranthene	12.60	202	449314	20.871	ng	99
78) Pyrene	12.83	202	379237	16.293	ng	98
81) Benzo(a)anthracene	14.02	228	212866	10.786	ng	98
83) Chrysene	14.06	228	167635	8.749	ng	98
86) Indeno(1,2,3-cd)pyrene	16.98	276	89853	5.583	ng	98
88) Benzo(b)fluoranthene	15.07	252	210047m	10.389	ng	
89) Benzo(k)fluoranthene	15.10	252	69999m	3.555	ng	
90) Benzo(a)pyrene	15.44	252	152393	8.432	ng	# 90
92) Benzo(a,h,i)perylene	17.43	276	86530	5.632	ng	97

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

