

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080219\
 Data File : BF115957.D
 Acq On : 2 Aug 2019 20:04
 Operator : HP/JU
 Sample : K4130-14
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 1S-L-(0-30)-COMP

Manual Integrations
 APPROVED

mohammad
 8/8/2019 4:44:32 PM

Quant Time: Aug 03 01:18:21 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	126124	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	493526	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	268318	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	493054	20.00	ng	0.00
76) Chrysene-d12	14.02	240	400543	20.00	ng	0.00
87) Perylene-d12	15.49	264	379548	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	739079	98.10	ng	0.00
7) Phenol-d6	6.48	99	971807	102.24	ng	0.00
23) Nitrobenzene-d5	7.41	82	605133	70.78	ng	0.00
42) 2,4,6-Tribromophenol	10.68	330	264356	101.62	ng	0.00
45) 2-Fluorobiphenyl	9.21	172	1033223	72.13	ng	0.00
79) Terphenyl-d14	12.96	244	1222796	64.33	ng	0.00
Target Compounds						
10) Phenol	6.49	94	53451	4.775	ng	91
50) Dimethylphthalate	9.60	163	87471	4.788	ng	99
71) Phenanthrene	11.40	178	90585	3.594	ng	97
75) Fluoranthene	12.59	202	243240	9.218	ng	98
78) Pyrene	12.82	202	211783	7.014	ng	98
81) Benzo(a)anthracene	14.01	228	116106	4.535	ng	99
83) Chrysene	14.05	228	88977	3.580	ng	98
86) Indeno(1,2,3-cd)pyrene	16.97	276	41843	2.004	ng	97
88) Benzo(b)fluoranthene	15.06	252	102083m	4.652	ng	
90) Benzo(a)pyrene	15.43	252	72102	3.675	ng	98
92) Benzo(g,h,i)perylene	17.41	276	40558	2.432	ng	97

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

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