

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081219\
 Data File : BF116213.D
 Acq On : 13 Aug 2019 2:23
 Operator : HP/JU
 Sample : K4292-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-COMP-WS

Manual Integrations
 APPROVED

mohammad
 8/14/2019 2:53:57 AM

Quant Time: Aug 13 03:58:06 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 11:13:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	123889	20.00	ng	-0.01
21) Naphthalene-d8	8.12	136	489141	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	248753	20.00	ng	-0.02
64) Phenanthrene-d10	11.36	188	394105	20.00	ng	-0.01
76) Chrysene-d12	14.00	240	276335	20.00	ng	-0.01
87) Perylene-d12	15.47	264	294353	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	728803	98.48	ng	-0.01
7) Phenol-d6	6.47	99	955245	102.31	ng	-0.01
23) Nitrobenzene-d5	7.40	82	601056	70.93	ng	-0.02
42) 2,4,6-Tribromophenol	10.66	330	227089	94.16	ng	-0.01
45) 2-Fluorobiphenyl	9.19	172	1079261	81.27	ng	-0.01
79) Terphenyl-d14	12.94	244	922826	70.37	ng	-0.01
Target Compounds						
10) Phenol	6.49	94	31540	2.868	ng	Qvalue 96
50) Dimethylphthalate	9.58	163	129146	7.625	ng	99
71) Phenanthrene	11.38	178	85655	4.251	ng	97
75) Fluoranthene	12.57	202	204525	9.697	ng	99
78) Pyrene	12.80	202	164650	7.904	ng	99
81) Benzo(a)anthracene	13.99	228	71521	4.049	ng	98
83) Chrysene	14.03	228	102249	5.963	ng	98
86) Indeno(1,2,3-cd)pyrene	16.96	276	53095	3.687	ng	96
88) Benzo(b)fluoranthene	15.04	252	154622m	9.085	ng	
89) Benzo(k)fluoranthene	15.07	252	46498m	2.805	ng	
90) Benzo(a)pyrene	15.41	252	77747	5.110	ng	# 94
92) Benzo(a,h,i)perylene	17.40	276	48367	3.740	ng	94

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

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