

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF052119\
 Data File : BF114470.D
 Acq On : 21 May 2019 21:35
 Operator : JU/SJ
 Sample : K2900-02
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P-2(0-10)

Manual Integrations
 APPROVED

Sohil
 5/23/2019 8:15:52 AM

Quant Time: May 22 07:42:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 20 04:47:40 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	205690	20.00	ng	-0.01
21) Naphthalene-d8	8.11	136	797557	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	360681	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	628490	20.00	ng	0.00
76) Chrysene-d12	14.00	240	411572	20.00	ng	0.00
87) Perylene-d12	15.47	264	366052	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	1134205	88.74	ng	0.00
7) Phenol-d6	6.47	99	1516855	100.34	ng	-0.01
23) Nitrobenzene-d5	7.39	82	909866	64.02	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	313041	83.95	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	1588759	66.63	ng	0.00
79) Terphenyl-d14	12.94	244	1197429	59.98	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.58	163	246274	9.301	ng	Qvalue 100
71) Phenanthrene	11.38	178	163564	5.349	ng	97
75) Fluoranthene	12.57	202	236618	7.186	ng	98
78) Pyrene	12.80	202	319295	9.644	ng	98
81) Benzo(a)anthracene	13.99	228	197082	7.403	ng	99
83) Chrysene	14.03	228	201374	7.717	ng	97
86) Indeno(1,2,3-cd)pyrene	16.95	276	96605	4.189	ng	97
88) Benzo(b)fluoranthene	15.04	252	274022m	11.685	ng	
89) Benzo(k)fluoranthene	15.07	252	69072m	3.206	ng	
90) Benzo(a)pyrene	15.41	252	170869	8.279	ng	100
92) Benzo(g,h,i)perylene	17.40	276	100546	5.850	ng	98

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

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