

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082819\
 Data File : BF116634.D
 Acq On : 29 Aug 2019 00:07
 Operator : HP/JU
 Sample : K4574-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S701A-N-0.0-0.5-082719

Manual Integrations
 APPROVED

Jagrut
 8/29/2019 7:47:23 PM

Quant Time: Aug 29 04:13:11 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 28 13:34:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	86445	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	256354	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	168101	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	228241	20.00	ng	0.00
76) Chrysene-d12	14.01	240	100818	20.00	ng	0.00
87) Perylene-d12	15.47	264	115236	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	589517	117.79	ng	0.00
7) Phenol-d6	6.49	99	756160	111.44	ng	0.00
23) Nitrobenzene-d5	7.42	82	368376	82.22	ng	0.00
42) 2,4,6-Tribromophenol	10.68	330	166599	104.15	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	656079	64.95	ng	0.00
79) Terphenyl-d14	12.96	244	571913	92.26	ng	0.00
Target Compounds						
10) Phenol	6.50	94	46965	6.310	ng	93
50) Dimethylphthalate	9.60	163	61908	5.478	ng	98
71) Phenanthrene	11.40	178	66119	5.393	ng	98
74) Di-n-butylphthalate	11.94	149	130094	10.760	ng	99
75) Fluoranthene	12.59	202	204305	18.582	ng	99
78) Pyrene	12.81	202	169166	18.248	ng	97
81) Benzo(a)anthracene	14.00	228	55489	8.255	ng	98
83) Chrysene	14.03	228	61470	9.342	ng	98
86) Indeno(1,2,3-cd)pyrene	16.91	276	30906	9.472	ng	96
88) Benzo(b)fluoranthene	15.04	252	88577m	12.517	ng	
89) Benzo(k)fluoranthene	15.07	252	29423m	4.374	ng	
90) Benzo(a)pyrene	15.40	252	56879	9.006	ng	96
92) Benzo(a,h,i)perylene	17.35	276	27385	4.469	ng	97

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

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