

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF053019\
 Data File : BF114671.D
 Acq On : 30 May 2019 20:22
 Operator : HP/JU
 Sample : K3065-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-16-COMP

Manual Integrations
 APPROVED

mohammad
 6/3/2019 8:47:18 AM

Quant Time: May 31 06:58:48 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 29 19:08:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	390841	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	1503030	20.00	ng	0.00
39) Acenaphthene-d10	9.72	164	757786	20.00	ng	0.00
64) Phenanthrene-d10	11.19	188	1478944	20.00	ng	0.00
76) Chrysene-d12	13.81	240	956415	20.00	ng	0.00
87) Perylene-d12	15.20	264	1048003	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	909436	40.85	ng	0.00
7) Phenol-d6	6.33	99	2159265	80.49	ng	0.00
23) Nitrobenzene-d5	7.25	82	1452577	60.29	ng	0.00
42) 2,4,6-Tribromophenol	10.50	330	185652	25.71	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	2385345	53.13	ng	0.00
79) Terphenyl-d14	12.77	244	2378384	46.75	ng	0.00
Target Compounds						
10) Phenol	6.33	94	75200	2.321	ng	96
50) Dimethylphthalate	9.44	163	220692	4.179	ng	98
75) Fluoranthene	12.39	202	217142	2.780	ng	98
78) Pyrene	12.62	202	242933	3.003	ng	95
88) Benzo(b)fluoranthene	14.81	252	135997m	2.060	ng	

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

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