

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082819\
 Data File : BF116637.D
 Acq On : 29 Aug 2019 1:29
 Operator : HP/JU
 Sample : K4576-02
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RBR-200034-RBR-200036

Manual Integrations
 APPROVED

Jagrut
 8/29/2019 7:47:30 PM

Quant Time: Aug 29 04:26:43 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	69317	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	281274	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	200948	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	168263	20.00	ng	0.00
76) Chrysene-d12	14.00	240	100854	20.00	ng	0.00
87) Perylene-d12	15.46	264	113424	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	611528	152.38	ng	0.00
7) Phenol-d6	6.49	99	643571	118.28	ng	0.00
23) Nitrobenzene-d5	7.42	82	405408	82.47	ng	0.00
42) 2,4,6-Tribromophenol	10.68	330	132926	69.51	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	704712	58.36	ng	0.00
79) Terphenyl-d14	12.96	244	423226	68.25	ng	0.00
Target Compounds						
10) Phenol	6.50	94	32168	5.390	ng	98
50) Dimethylphthalate	9.60	163	96944	7.176	ng	99
71) Phenanthrene	11.40	178	19864	2.198	ng	98
75) Fluoranthene	12.59	202	38127	4.704	ng	96
78) Pyrene	12.81	202	29269	3.156	ng	96
83) Chrysene	14.03	228	13286	2.018	ng	95
86) Indeno(1,2,3-cd)pyrene	16.90	276	4385	4.652	ng	95
88) Benzo(b)fluoranthene	15.03	252	15165m	2.177	ng	

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

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