

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081019\
 Data File : BF116173.D
 Acq On : 11 Aug 2019 4:05
 Operator : HP/JU
 Sample : K4293-01
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PE-1

Manual Integrations
 APPROVED

mohammad
 8/14/2019 2:55:08 AM

Quant Time: Aug 12 04:30:54 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	6.84	152	152676	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	585694	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	281471	20.00	ng	-0.02
64) Phenanthrene-d10	11.36	188	414745	20.00	ng	-0.01
76) Chrysene-d12	14.00	240	286140	20.00	ng	-0.01
87) Perylene-d12	15.47	264	276883	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	779193	85.44	ng	0.00
7) Phenol-d6	6.47	99	998329	86.76	ng	-0.01
23) Nitrobenzene-d5	7.40	82	652899	64.35	ng	-0.02
42) 2,4,6-Tribromophenol	10.66	330	185663	68.03	ng	-0.01
45) 2-Fluorobiphenyl	9.19	172	1135537	75.57	ng	-0.01
79) Terphenyl-d14	12.94	244	925938	68.19	ng	-0.01
Target Compounds						
10) Phenol	6.49	94	30399	2.243	ng	95
50) Dimethylphthalate	9.59	163	106407	5.552	ng	99
75) Fluoranthene	12.57	202	88624	3.993	ng	99
78) Pyrene	12.80	202	85922	3.983	ng	97
81) Benzo(a)anthracene	13.99	228	47352	2.589	ng	97
83) Chrysene	14.03	228	46260	2.605	ng	98
88) Benzo(b)fluoranthene	15.04	252	56809m	3.548	ng	
90) Benzo(a)pyrene	15.41	252	44628	3.118	ng	99

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

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