

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122019\
 Data File : BP001353.D
 Acq On : 20 Dec 2019 20:14
 Operator : JU
 Sample : K6107-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 HR-06-121819

Manual Integrations
 APPROVED

mohammad
 12/23/2019 11:56:18 AM

Quant Time: Dec 21 03:48:26 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 20 03:20:39 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.30	152	43200	20.00	ng	0.00
21) Naphthalene-d8	10.33	136	151173	20.00	ng	0.00
39) Acenaphthene-d10	13.20	164	85850	20.00	ng	0.00
64) Phenanthrene-d10	15.60	188	163222	20.00	ng	0.00
76) Chrysene-d12	19.25	240	137136	20.00	ng	0.00
87) Perylene-d12	21.02	264	149836	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.21	112	379337	141.92	ng	0.01
7) Phenol-d6	7.76	99	519088	148.83	ng	0.00
23) Nitrobenzene-d5	9.19	82	353691	89.85	ng	0.00
42) 2,4,6-Tribromophenol	14.50	330	142820	133.06	ng	0.00
45) 2-Fluorobiphenyl	12.12	172	557957	82.18	ng	0.00
79) Terphenyl-d14	17.89	244	660671	82.51	ng	0.00
Target Compounds						
10) Phenol	7.77	94	22154	5.559	ng	84
37) 2-Methylnaphthalene	11.50	142	12673	2.225	ng	# 80
50) Dimethylphthalate	12.79	163	51879	7.366	ng	# 98
71) Phenanthrene	15.63	178	27354	2.937	ng	99
75) Fluoranthene	17.35	202	51298	4.926	ng	98
78) Pyrene	17.65	202	46550	4.372	ng	# 95
81) Benzo(a)anthracene	19.23	228	25895	2.591	ng	95
83) Chrysene	19.27	228	22793m	2.363	ng	
88) Benzo(b)fluoranthene	20.53	252	26486m	2.692	ng	
90) Benzo(a)pyrene	20.94	252	18971	2.116	ng	# 89

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

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