

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
 Data File : BP001399.D
 Acq On : 24 Dec 2019 21:34
 Operator : JU
 Sample : K6413-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 K005-AA-WC-2

Manual Integrations
 APPROVED

mohammad
 12/26/2019 4:01:37 PM

Quant Time: Dec 25 05:40:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 16:04:49 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.28	152	51769	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	192634	20.00	ng	0.00
39) Acenaphthene-d10	13.18	164	116318	20.00	ng	0.00
64) Phenanthrene-d10	15.57	188	219570	20.00	ng	0.00
76) Chrysene-d12	19.22	240	143795	20.00	ng	0.00
87) Perylene-d12	20.99	264	142472	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.20	112	266058	83.06	ng	0.00
7) Phenol-d6	7.75	99	375068	89.74	ng	0.00
23) Nitrobenzene-d5	9.16	82	257886	51.41	ng	-0.01
42) 2,4,6-Tribromophenol	14.47	330	118128	81.23	ng	0.00
45) 2-Fluorobiphenyl	12.10	172	485069	52.73	ng	0.00
79) Terphenyl-d14	17.86	244	541296	64.47	ng	0.00
Target Compounds						
10) Phenol	7.76	94	12770	2.674	ng	84
50) Dimethylphthalate	12.76	163	32132	3.367	ng	# 96
71) Phenanthrene	15.61	178	40516	3.234	ng	95
75) Fluoranthene	17.32	202	62258	4.445	ng	99
78) Pyrene	17.63	202	55811	4.999	ng	# 95
81) Benzo(a)anthracene	19.21	228	26335	2.513	ng	95
83) Chrysene	19.25	228	24011	2.374	ng	98
88) Benzo(b)fluoranthene	20.50	252	27661m	2.956	ng	
90) Benzo(a)pyrene	20.92	252	19907	2.336	ng	# 98

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

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