

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122319\
 Data File : BP001380.D
 Acq On : 23 Dec 2019 21:37
 Operator : JU
 Sample : K6117-07
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TR-04-121919

Manual Integrations
 APPROVED

mohammad
 12/24/2019 3:05:39 PM

Quant Time: Dec 24 06:26:21 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	58015	20.00	ng	0.00
21) Naphthalene-d8	10.34	136	204955	20.00	ng	0.00
39) Acenaphthene-d10	13.20	164	118665	20.00	ng	0.00
64) Phenanthrene-d10	15.60	188	207121	20.00	ng	0.00
76) Chrysene-d12	19.26	240	142091	20.00	ng	0.01
87) Perylene-d12	21.03	264	151854	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	6.22	112	335767	93.54	ng	0.02
7) Phenol-d6	7.77	99	466339	99.56	ng	0.02
23) Nitrobenzene-d5	9.19	82	325807	61.04	ng	0.00
42) 2,4,6-Tribromophenol	14.50	330	126228	85.08	ng	0.01
45) 2-Fluorobiphenyl	12.12	172	595265	63.43	ng	0.00
79) Terphenyl-d14	17.90	244	568189	68.49	ng	0.01
Target Compounds						
10) Phenol	7.78	94	14803	2.766	ng	83
49) Acenaphthylene	12.97	152	66803	5.693	ng	98
50) Dimethylphthalate	12.79	163	24776	2.545	ng	# 99
71) Phenanthrene	15.64	178	27389	2.318	ng	99
72) Anthracene	15.72	178	33606	2.873	ng	99
75) Fluoranthene	17.36	202	76764	5.810	ng	99
78) Pyrene	17.66	202	134075	12.153	ng	97
81) Benzo(a)anthracene	19.25	228	67642	6.533	ng	# 78
83) Chrysene	19.29	228	71406	7.145	ng	# 93
86) Indeno(1,2,3-cd)pyrene	22.67	276	42053	3.896	ng	99
88) Benzo(b)fluoranthene	20.55	252	74684m	7.489	ng	
89) Benzo(k)fluoranthene	20.57	252	19169m	2.018	ng	
90) Benzo(a)pyrene	20.96	252	63634	7.004	ng	# 96
92) Benzo(a,h,i)perylene	23.17	276	52031	6.130	ng	100

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

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