

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100820\
 Data File : BP003580.D
 Acq On : 08 Oct 2020 20:05
 Operator : CG/JU
 Sample : L4291-11
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PAV-B4-100620-DP

Manual Integrations
 APPROVED

mohammad
 10/12/2020 11:32:29 AM

Quant Time: Oct 09 01:59:35 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 08 15:55:58 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.481	152	52147	20.00	ng	0.00
21) Naphthalene-d8	10.257	136	197933	20.00	ng	0.00
39) Acenaphthene-d10	14.139	164	117313	20.00	ng	0.00
64) Phenanthrene-d10	16.898	188	257780	20.00	ng	0.00
76) Chrysene-d12	21.010	240	298693	20.00	ng	# 0.00
86) Perylene-d12	23.174	264	346635	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.099	112	212899	71.29	ng	0.00
7) Phenol-d6	6.693	99	249208	62.60	ng	0.00
23) Nitrobenzene-d5	8.634	82	184908	54.88	ng	0.00
42) 2,4,6-Tribromophenol	15.651	330	116429	65.07	ng	0.00
45) 2-Fluorobiphenyl	12.751	172	406489	50.68	ng	0.00
79) Terphenyl-d14	19.486	244	636538	44.40	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	13.610	163	46340	4.95	ng	99
71) Phenanthrene	16.939	178	156381	11.15	ng	100
72) Anthracene	17.039	178	39423	2.85	ng	99
75) Fluoranthene	18.933	202	361019	20.47	ng	98
78) Pyrene	19.286	202	361986	19.43	ng	100
81) Benzo(a)anthracene	20.998	228	260485	13.49	ng	99
83) Chrysene	21.045	228	230819	12.44	ng	97
87) Indeno(1,2,3-cd)pyrene	25.368	276	215572	8.20	ng	# 92
88) Benzo(b)fluoranthene	22.533	252	372078	17.00	ng	98
89) Benzo(k)fluoranthene	22.568	252	126809m	6.04	ng	
90) Benzo(a)pyrene	23.080	252	311191	15.08	ng	95
91) Dibenzo(a,h)anthracene	25.362	278	65566m	3.02	ng	
92) Benzo(g,h,i)perylene	26.039	276	210796	9.74	ng	# 94

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

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