

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100820\
 Data File : BP003576.D
 Acq On : 08 Oct 2020 17:49
 Operator : CG/JU
 Sample : L4292-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PAV-B13-100620-0-10

Manual Integrations
 APPROVED

mohammad
 10/12/2020 11:32:13 AM

Quant Time: Oct 08 18:38:18 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	59168	20.00	ng	0.00
21) Naphthalene-d8	10.257	136	208516	20.00	ng	# 0.00
39) Acenaphthene-d10	14.139	164	115378	20.00	ng	0.00
64) Phenanthrene-d10	16.898	188	244778	20.00	ng	0.00
76) Chrysene-d12	21.010	240	289882	20.00	ng	0.00
86) Perylene-d12	23.174	264	346119	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.099	112	311459	91.92	ng	0.00
7) Phenol-d6	6.693	99	350143	77.52	ng	0.00
23) Nitrobenzene-d5	8.634	82	258032	72.70	ng	0.00
42) 2,4,6-Tribromophenol	15.645	330	153321	87.12	ng	0.00
45) 2-Fluorobiphenyl	12.751	172	528170	66.95	ng	0.00
79) Terphenyl-d14	19.486	244	791040	56.86	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	13.610	163	50217	5.45	ng	99
71) Phenanthrene	16.939	178	62500	4.69	ng	99
75) Fluoranthene	18.933	202	147611	8.81	ng	99
78) Pyrene	19.280	202	165211	9.14	ng	99
81) Benzo(a)anthracene	20.992	228	102039	5.45	ng	# 93
83) Chrysene	21.045	228	101213	5.62	ng	# 93
87) Indeno(1,2,3-cd)pyrene	25.380	276	98779	3.77	ng	# 90
88) Benzo(b)fluoranthene	22.533	252	161388	7.38	ng	# 93
89) Benzo(k)fluoranthene	22.568	252	48788m	2.33	ng	# 93
90) Benzo(a)pyrene	23.086	252	139752	6.78	ng	# 93
92) Benzo(g,h,i)perylene	26.051	276	111478	5.16	ng	94

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

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